

Z I N C

General analytical conditions

STOCK SOLUTION

Dissolve 1.000 g metallic Zn in about 3 ml conc. HC 1 and make up to 1 l.
This solution contains 1000 mg/l Zn.

INTS ON ANALYSIS

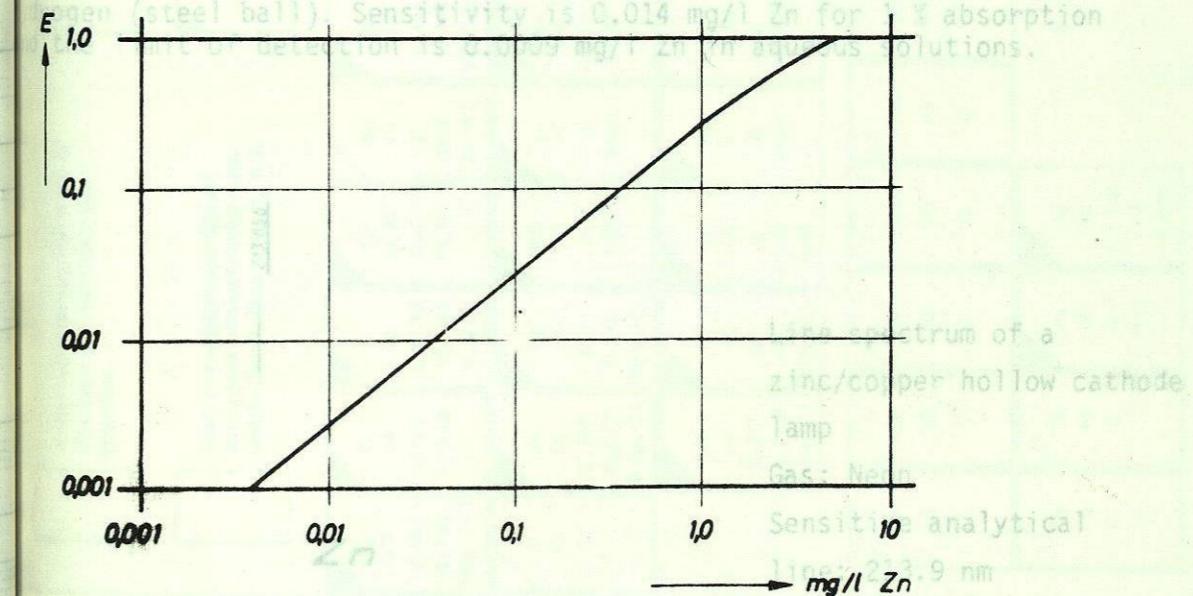
INSTRUMENT SETTING

Wavelength	427.7	213.9 nm
Barrier filter	I	
Slit	0.1 mm	0.3 nm
Burner		
Burner height	universal	
Support gas	15.5	
Fuel gas	8 - 9 *	air
Lamp current	15 mA	acetylene

* Set to minimum flame absorption with blank solution.

TYPICAL CALIBRATION CURVE with the instrument setting given

Maximum reading stability and limit of detection is obtained when using universal burner with 16.5 scale divisions air and 5-6 scale divisions propane (steel ball). Sensitivity is 0.014 mg/l Zn for 1 % absorption



SENSITIVITY 0.012 mg/l Zn for 1 % absorption in aqueous solution.

LIMIT OF DETECTION 0.002 mg/l Zn in aqueous solution.

ZINC

General analytical conditions

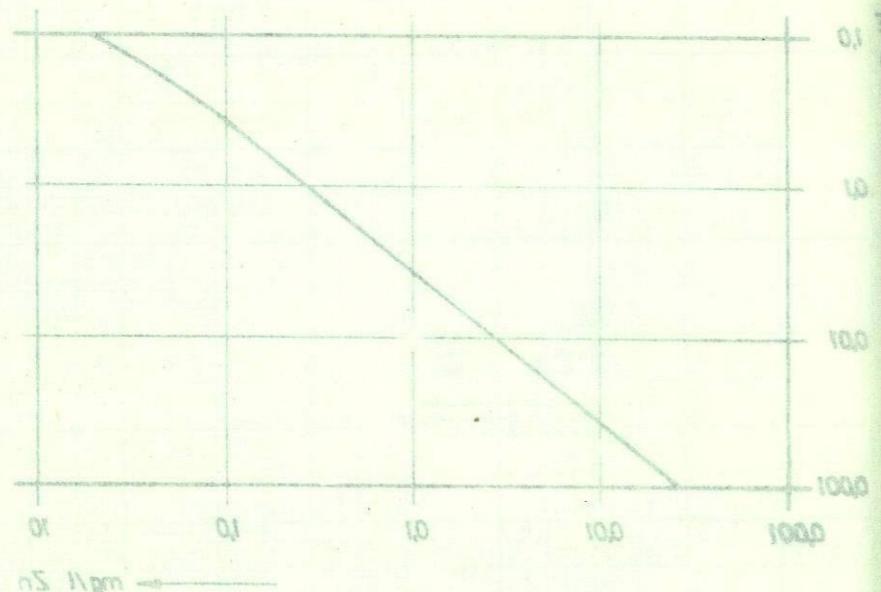
Assay 1.000 g metallic Zn in spot 3 ml conc. HCl and make up to 1 l.

Solutions 1000 mg/l Zn.

mm 0.3	AS 5.0	Magnetruck Burner filter Slit
mm 8.0	I 1.0	
	Universal gas - 8	
	12.5 * 8	
acetylene burner	12.5	Support gas
hollow	12 Am	Fuel gas
		Lamp current

Set to minimum flame absorption with blank solution.

PIRICAL CALIBRATION CURVE with the instrument setting given



Set to 1% absorption in aqueous solution.

LIMIT OF DETECTION 0.005 mg/l Zn in aqueous solution.

BLACK SOLUTION

INTERFERENCES

So far no interferences have been observed in the air/acetylene flame.

INSTRUMENT SETTING

INTS ON ANALYSIS

Slit width:

Slit width can be increased to 0.5 mm if Cu-free hollow cathode lamp is used.

Lamp current:

With increasing current the sensitivity of the measuring set-up is reduced.
ideal amperage: 7 - 10 mA.

Air/propane flame:

Sensitivity is by about 40 % higher than in the air/acetylene flame with propane burner and with 15.5 scale divisions air and 4-5 scale divisions propane (set to minimum intrinsic absorption of flame). But the worse reading stability does not improve the detection limit.

Air/hydrogen flame:

Maximum reading stability and limit of detection is obtained when using universal burner with 16.5 scale divisions air and 5-6 scale divisions hydrogen (steel ball). Sensitivity is 0.014 mg/l Zn for 1 % absorption and the limit of detection is 0.0009 mg/l Zn in aqueous solutions.

Atomic-absorption absorption setting on the signal not required to obtain flame.

Weak or negligible absorption on inter-base anion-exchange resin.

Weight fraction 213.856 < 5

Li +1 C	Na +2 C	Br -1 C	Al +3 C	Tl +3 C
K +1 C	Mg +2 C	Ca +2 C	Si +4 C	Y +3 C
Rb +1 C	Ca +2 C	Al +3 C	Sn +4 C	Ge +4 C
Cs +1 C	Al +3 C	Si +4 C	Pb +4 C	Pr +4 C

Zn

Line spectrum of a zinc/copper hollow cathode lamp
Gas: Neon
Sensitive analytical line: 213.9 nm

Anion-exchange behavior in chloride systems (0-12 Nolar)

Fig. 1

Fig. 1

Anion-exchange behavior in chloride systems (0.12 Molar)

ANION PREPARATION			TABLE IV. RETENTION TIMES ON PORAPAK POROUS POLYMERS (MINUTES)											
L	B	C	S	P	Br	Ge	As	Se	Te	At	Lu	S	N	T
L	+1	C	Si	C	A > 5	C, 0.8	A > 5	A > 3	A > 4	A	Yb	+3	C	0.8
Mg	+2	C	Co	+2	C, 0.1	A, 4-12	C, 0.1	Cd	+2	A > 5	Tm	+3	C	0.8
Na	+1	C	Fe	+3	C, 0-1	A, 1-12	C, 0-1	In	+3	C	Er	+3	C	0.8
K	+1	C	Mn	+2	C, 0-4	A, 4-12	C, 0-4	Sn	+3	A > 5	Dy	+3	C	0.8
Rb	+1	C	Cr	+4	A, 11-12	+5 Red.	A, 4-12	Ge	+4	A > 3	Tb	+3	C	0.8
Sc	+3	C	V	+4	A, 11-12	+5 Red.	A, 9-12	As	+3	A > 5	Ho	+3	C	0.8
Y	+3	C	Ti	+4	A, 9-12	+5 Red.	A, 9-12	Se	+4	A > 3	Eu	+3	C	0.8
Zr	+4	C	Sc	+3	C, 0-7	A, 7-12	Poly. < 1	Pd	+2	A > 6M	Pm	+3	C	0.8
Nb	+5	A > 12	V	+4	A, 11-12	+5 Red.	A, 9-12	Ag	+1	Inert	Nd	+3	C	0.8
Mo	+6	A	Cr	+4	A, 11-12	+5 Red.	A, 9-12	Rh	+4	A > 12	Pr	+4	C	0.8
Tc	+7	A	Mn	+2	C, 0-1	A, 6-12	C, 0-1	Ru	+4	A > 1 M	Ta	+5	C	0.8
Ru	+4	A > 1 M	Co	+2	C, 0-6	A, 6-12	C, 0-6	Tc	+3	A > 6M	Hf	+4	C	0.8
Tc	+7	A	Fe	+3	C, 0-1	A, 6-12	C, 0-1	Ir	+4	A > 0.1 M	Os	+3, +4	C	0.8
Re	+7	A	Mn	+2	C, 0-1	A, 6-12	C, 0-1	Pt	+4	A > 0.1 M	Ta	+5	C	0.8
W	+6	A	Sc	+3	C, 0-7	A, 7-12	Poly. < 1	Pt	+4	A > 0.1 M	Hf	+4	C	0.8
Poly.	Hyd.		Sc	+3	C, 0-7	A, 7-12	Poly. < 1	Re	+7	A	Os	+3, +4	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1	Re	+7	A	Ta	+5	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1	W	+6	A	Hf	+4	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1	Poly.	Hyd.		Os	+3, +4	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1				Ta	+5	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1				Hf	+4	C	0.8
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			Sc	+3	C, 0-7	A, 7-12	Poly. < 1				Ta	+5	C	0.8
			Sc	+3	C, 0-7	A, 7-12	Poly. < 1				Hf	+4	C	0

MN PREPARATION

RAPAK beads do not crush or
easily and may be packed
into GC columns. For best re-
sults, the packing material should
be added to the column through a
hopper in small amounts — a fraction
of a cubic centimeter. This is
enough material to fill 5 to 7.5 cm
(3 in.) in column length at a
time. After each addition, the col-
umn should be vibrated or gently
tapped to insure the column is
packed tightly and has a low void
volume. The efficiency of the col-
umn is directly dependent upon how
well the column is packed.

column should be conditioned carrier flow for at least two hours at about 25 to 50°C below the temperature limit of the PORA-1. While the column is conditioned, the effluent end should not be directed to the detector. For very efficient columns more methanol should be added to the effluent of the column at the end of the conditioning period to fill any voids. The column should be further conditioned for an additional half hour. At highest sensitivity, the column should be conditioned overnight. Porous polymers will discolor on heating, but this does not affect the separation.

TABLE III. PHYSICAL PROPERTIES FOR PAPAK POROUS POLYMERS

Surface (m ² /gm)	Density (g/cm ³)	Maximum Temperature (°C)
100-200	0.27	250
500-600	0.34	250
450-600	0.30	250
300-450	0.35	250
225-350	0.38	190
250-350	0.43	190

TABLE IV. RETENTION TIMES
ON PORAPAK POROUS POLYMERS (MINUTES)

COMPOUND		P	Q	R	S	N	T
ACIDS	Acetic Acid	0.39	1.08	2.14	3.68	4.18	5.21
	Propionic Acid	0.65	2.42	4.75	9.05	9.11	10.68
	iso-Butyric Acid	0.91	4.43	8.28	16.41	16.32	18.13
	Butyric Acid	1.05	5.22	9.87	19.71	19.80	22.30
	iso-Valeric Acid	1.52	9.69	16.85	36.81	36.79	39.50
	Valeric Acid	1.89	12.12	21.70	45.70	45.30	48.90
ALCOHOLS	Methanol	0.24	0.35	0.39	0.48	0.61	0.69
	Ethanol	0.30	0.60	0.63	0.76	1.21	1.20
	iso-Propyl alcohol	0.37	0.91	0.89	1.03	1.79	1.90
	tert-Butyl alcohol	0.38	1.40	1.35	1.69	2.56	2.67
	n-Propyl alcohol	0.43	1.22	1.24	1.53	2.52	2.60
	sec-Butyl alcohol	0.50	1.93	1.85	2.40	3.69	3.81
	iso-Butyl alcohol	0.61	2.18	2.16	2.72	4.63	4.74
	tert-Pentyl alcohol	0.62	3.17	2.90	3.78	5.87	5.84
	n-Butyl alcohol	0.69	2.58	2.68	3.39	5.78	5.71
	iso-Pentyl alcohol	1.07	4.87	4.75	6.14	10.82	10.76
	n-Pentyl alcohol	1.10	5.63	5.67	7.04	12.88	12.40
ALDEHYDES	Acetaldehyde	0.24	0.44	0.39	0.49	0.65	0.71
	Propionaldehyde	0.35	0.87	0.79	0.97	1.44	1.46
	Acrolein	0.35	0.81	0.77	0.91	1.44	1.52
	iso-Butyraldehyde	0.46	1.51	1.37	1.67	2.54	2.50
	Butyraldehyde	0.54	1.77	1.62	1.96	3.07	3.04
ALKANES	Pentane	0.34	1.24	1.02	1.19	1.39	1.10
	Hexane	0.48	2.56	2.03	2.46	3.00	2.20
	Heptane	0.73	5.26	4.27	5.03	6.22	4.39
	Octane	1.14	10.99	8.54	10.36	13.28	8.94
ARYLS	Benzene	0.87	2.76	2.56	3.00	4.00	3.74
	Toluene	1.33	5.72	5.31	6.41	8.54	7.61
	Ethyl Benzene	2.09	11.60	10.69	12.98	17.58	15.05
	o-Xylene	2.52	13.48	12.53	15.32	20.79	18.05
CHLOROALKYLS	Methylene Chloride	0.41	0.98	0.94	1.09	1.48	1.67
	Chloroform	0.59	1.89	1.76	2.15	2.91	3.04
	Carbon Tetrachloride	0.75	2.87	2.42	2.98	3.49	3.04
	1,2-Dichloroethane	0.78	2.28	2.16	2.56	3.84	4.09
	1,2-Dichloropropane	0.98	3.84	3.47	4.23	6.29	6.37
ESTERS	Methyl formate	0.27	0.51	0.53	0.57	0.91	0.94
	Methyl acetate	0.38	1.01	0.97	1.11	1.74	1.75
	Ethyl acetate	0.52	1.99	1.83	2.15	3.30	3.16
	Methyl propionate	0.57	1.99	1.83	2.15	3.30	3.16
	iso-Propyl acetate	0.65	3.18	2.75	3.38	5.22	4.89
	Propyl acetate	0.83	4.09	3.56	4.36	6.91	6.57
	Butyl acetate	1.39	8.78	7.35	9.44	14.60	13.88
ETHERS	Diethyl ether	0.33	1.41	0.98	1.27	1.45	1.28
	iso-Propyl ether	0.52	2.99	2.40	2.92	3.94	3.20
	Dibutyl ether	1.94	19.64	14.55	19.11	26.95	20.25
GLYCOLS	Ethylene glycol	0.94	2.39	3.36	4.19	9.34	12.43
	2,3-Butanediol	1.41	6.06	-	-	-	-
	1,3-Propanediol	1.83	6.06	-	-	-	-
	1,3-Butanediol	2.19	9.64	-	-	-	-
	1,4-Butanediol	3.38	13.97	-	-	-	-
	Diethylene glycol	4.21	16.95	-	-	-	-
GLYCEROL	Glycerol	4.92	19.15	-	-	-	-
KETONES	Acetone	0.33	0.89	0.81	0.96	1.59	1.66
	2-Butanone	0.53	1.82	1.59	1.95	3.22	3.17
	2-Pentanone	0.80	3.67	3.17	4.02	6.63	6.32
NITRILES	Acetonitrile	0.38	0.75	0.76	0.87	1.50	1.85
	Propionitrile	0.44	1.02	1.01	1.16	1.97	2.23
	Acrylonitrile	0.55	1.40	1.37	1.62	2.91	3.34
ADDITIONAL COMPOUNDS	Nitromethane	0.53	1.02	1.13	1.26	2.54	2.99
	Dioxane	0.98	3.35	2.84	3.39	5.28	5.91
	Pyridine	1.35	4.34	4.06	4.59	7.53	7.24
	Dimethyl sulfoxide	2.60	8.70	7.85	8.38	19.50	23.62
	Phenol	4.04	15.11	-	-	-	-
	Benzyl alcohol	6.73	30.62	-	-	-	-

Operating Conditions: 1 molar C_2H_2 at 23 mm Hg, stainless steel column, 80–100 mm tube, 268°C, 25–11 Hz, $\text{N}_2/\text{O}_2 = 5/1$.

TABLE I. RETENTION TIMES
ON PORAPAK POROUS POLYMERS (MINUTES)

T	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	25	26	27	28	29	30	31	32	33	34	35	36	37	38	39	40	41	42	43	44	45	46	47	48	49	50	51	52	53	54	55	56	57	58	59	60	61	62	63	64	65	66	67	68	69	70	71	72	73	74	75	76	77	78	79	80	81	82	83	84	85	86	87	88	89	90	91	92	93	94	95	96	97	98	99	100	101	102	103	104	105	106	107	108	109	110	111	112	113	114	115	116	117	118	119	120	121	122	123	124	125	126	127	128	129	130	131	132	133	134	135	136	137	138	139	140	141	142	143	144	145	146	147	148	149	150	151	152	153	154	155	156	157	158	159	160	161	162	163	164	165	166	167	168	169	170	171	172	173	174	175	176	177	178	179	180	181	182	183	184	185	186	187	188	189	190	191	192	193	194	195	196	197	198	199	200	201	202	203	204	205	206	207	208	209	210	211	212	213	214	215	216	217	218	219	220	221	222	223	224	225	226	227	228	229	230	231	232	233	234	235	236	237	238	239	240	241	242	243	244	245	246	247	248	249	250	251	252	253	254	255	256	257	258	259	260	261	262	263	264	265	266	267	268	269	270	271	272	273	274	275	276	277	278	279	280	281	282	283	284	285	286	287	288	289	290	291	292	293	294	295	296	297	298	299	300	301	302	303	304	305	306	307	308	309	310	311	312	313	314	315	316	317	318	319	320	321	322	323	324	325	326	327	328	329	330	331	332	333	334	335	336	337	338	339	340	341	342	343	344	345	346	347	348	349	350	351	352	353	354	355	356	357	358	359	360	361	362	363	364	365	366	367	368	369	370	371	372	373	374	375	376	377	378	379	380	381	382	383	384	385	386	387	388	389	390	391	392	393	394	395	396	397	398	399	400	401	402	403	404	405	406	407	408	409	410	411	412	413	414	415	416	417	418	419	420	421	422	423	424	425	426	427	428	429	430	431	432	433	434	435	436	437	438	439	440	441	442	443	444	445	446	447	448	449	450	451	452	453	454	455	456	457	458	459	460	461	462	463	464	465	466	467	468	469	470	471	472	473	474	475	476	477	478	479	480	481	482	483	484	485	486	487	488	489	490	491	492	493	494	495	496	497	498	499	500	501	502	503	504	505	506	507	508	509	510	511	512	513	514	515	516	517	518	519	520	521	522	523	524	525	526	527	528	529	530	531	532	533	534	535	536	537	538	539	540	541	542	543	544	545	546	547	548	549	550	551	552	553	554	555	556	557	558	559	560	561	562	563	564	565	566	567	568	569	570	571	572	573	574	575	576	577	578	579	580	581	582	583	584	585	586	587	588	589	590	591	592	593	594	595	596	597	598	599	600	601	602	603	604	605	606	607	608	609	610	611	612	613	614	615	616	617	618	619	620	621	622	623	624	625	626	627	628	629	630	631	632	633	634	635	636	637	638	639	640	641	642	643	644	645	646	647	648	649	650	651	652	653	654	655	656	657	658	659	660	661	662	663	664	665	666	667	668	669	670	671	672	673	674	675	676	677	678	679	680	681	682	683	684	685	686	687	688	689	690	691	692	693	694	695	696	697	698	699	700	701	702	703	704	705	706	707	708	709	710	711	712	713	714	715	716	717	718	719	720	721	722	723	724	725	726	727	728	729	730	731	732	733	734	735	736	737	738	739	740	741	742	743	744	745	746	747	748	749	750	751	752	753	754	755	756	757	758	759	760	761	762	763	764	765	766	767	768	769	770	771	772	773	774	775	776	777	778	779	770	771	772	773	774	775	776	777	778	779	780	781	782	783	784	785	786	787	788	789	790	791	792	793	794	795	796	797	798	799	800	801	802	803	804	805	806	807	808	809	810	811	812	813	814	815	816	817	818	819	820	821	822	823	824	825	826	827	828	829	830	831	832	833	834	835	836	837	838	839	830	831	832	833	834	835	836	837	838	839	840	841	842	843	844	845	846	847	848	849	850	851	852	853	854	855	856	857	858	859	850	851	852	853	854	855	856	857	858	859	860	861	862	863	864	865	866	867	868	869	860	861	862	863	864	865	866	867	868	869	870	871	872	873	874	875	876	877	878	879	870	871	872	873	874	875	876	877	878	879	880	881	882	883	884	885	886	887	888	889	880	881	882	883	884	885	886	887	888	889	890	891	892	893	894	895	896	897	898	899	890	891	892	893	894	895	896	897	898	899	900	901	902	903	904	905	906	907	908	909	900	901	902	903	904	905	906	907	908	909	910	911	912	913	914	915	916	917	918	919	910	

BRESCENTS PROPERTIES OF THE L-TAKTAROL AMINO ACIDS

DETERMINACION DE HIERRO

1,10 Fenantrolina: 0.2 gramos en 100 ml de agua caliente.

Clorhidrato de Hidroxilamina: 5 gramos en 100 ml de agua destilada.

Acetato de Sodio 1M : Disolver 14 gramos en 100 ml de agua destilada.

Solución Patrón de Hierro (100 ppm): Disolver 0.702 gramos de $\text{FeSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4 \cdot 6\text{H}_2\text{O}$
 (P.M.=392.14) en 100 ml de Ácido Sulfúrico -

(P.M.=392.14) en 100 ml de Acido Sulfurico -

0.1M y luego aforar a un litro con el mismo ácido.

ELECTROFORESIS

Aminoácidos: Acido aspártico, Histidina, Lisina y una mezcla de los tres en Amortiguador Tris-Acetato 0.07M pH 7.6 que contenga 10 g/l de

Amortiguador Tris-Acetato 0.07M pH 7.6 que contenga 10 g/l de Glucosa.

Minhidrina: Disolver 0.2 gramos en 100 ml de acetona.

Anilina-Difenilamina-Acido Fosfórico (5:5:1): Mezclar en las proporciones

indicadas las siguientes soluciones:

Anilina 10 g/l en acetona, Difenilamina -

10 g/l en acetona y Ácido Fosfórico 85%.

