

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
 Data File : BM014488.D
 Acq On : 06 Mar 2018 00:23
 Operator : SJ/JU
 Sample : J1795-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SVOC-GPC2-BLANK

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	64	66	67	rBV2	11267	10577	1.11%	0.198%
2	4.987	339	340	344	rBV3	7932	10751	1.13%	0.201%
3	7.510	763	769	781	rBV	256073	512754	53.98%	9.583%
4	10.286	1233	1241	1252	rBV	272342	625427	65.84%	11.688%
5	14.174	1895	1902	1912	rBV3	388013	773056	81.38%	14.447%
6	16.733	2334	2337	2344	rVB8	11111	22171	2.33%	0.414%
7	16.945	2367	2373	2393	rBV	454078	949924	100.00%	17.753%
8	17.109	2398	2401	2403	rBV4	10755	11866	1.25%	0.222%
9	17.209	2417	2418	2423	rBV4	9084	12302	1.30%	0.230%
10	17.956	2544	2545	2550	rBV5	9802	12781	1.35%	0.239%
11	18.380	2614	2617	2618	rBV3	10521	10387	1.09%	0.194%
12	18.809	2689	2690	2693	rBV3	11889	11238	1.18%	0.210%
13	19.397	2787	2790	2794	rBV5	54590	53414	5.62%	0.998%
14	20.562	2984	2988	2992	rBV	632955	708868	74.62%	13.248%
15	21.156	3085	3089	3096	rBV2	615108	853202	89.82%	15.945%
16	23.338	3455	3460	3468	rVB	395223	772184	81.29%	14.431%

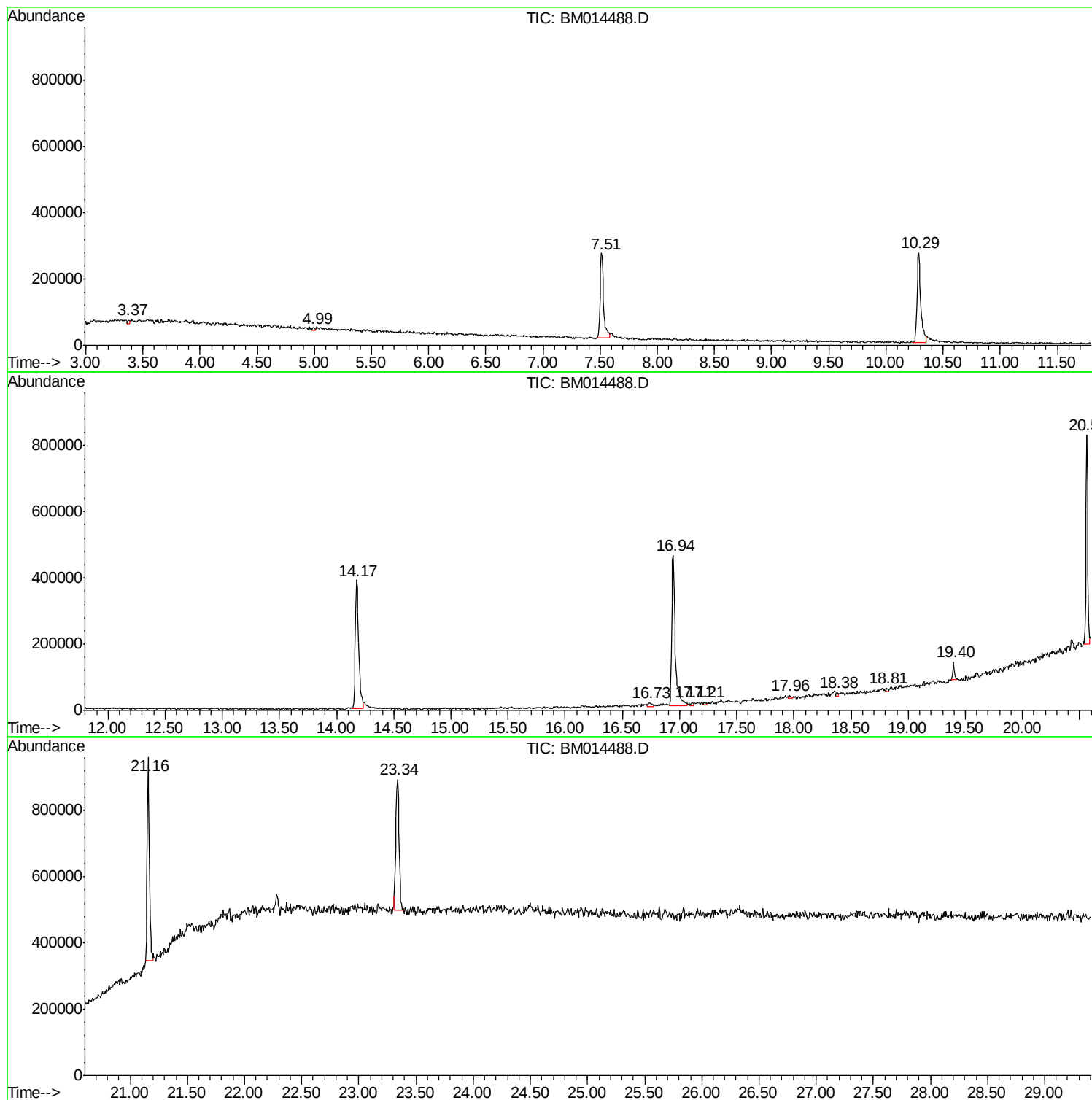
Sum of corrected areas: 5350902

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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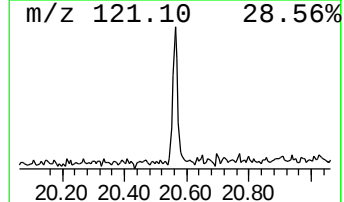
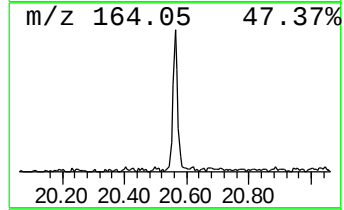
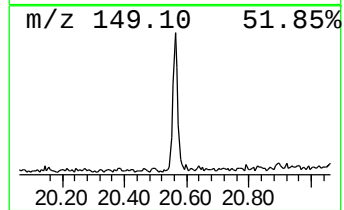
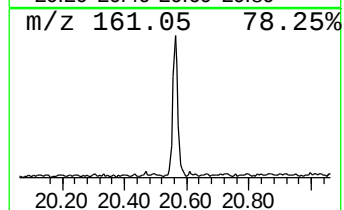
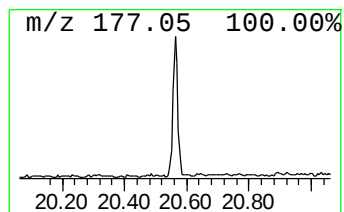
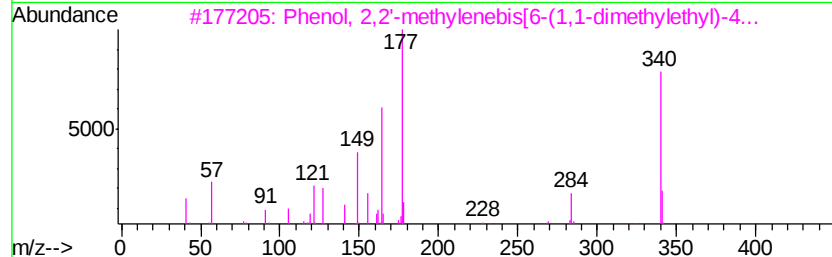
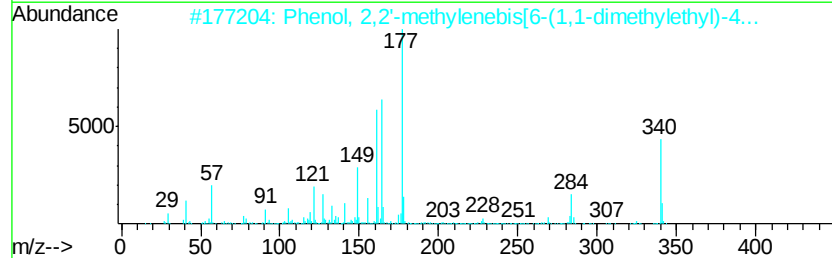
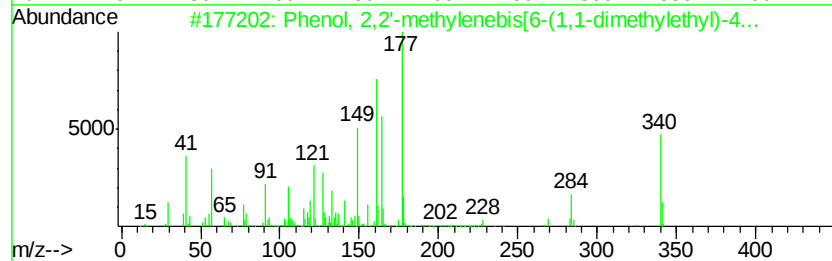
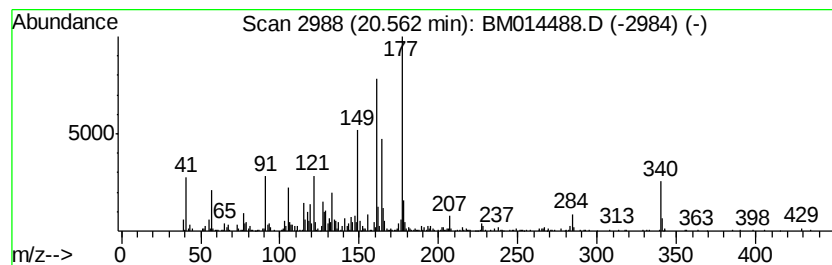
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenol, 2,2'-methylenebis[6... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.56	16.62 ng/ul	708868	Chrysene-d12	21.16		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,2'-methylenebis[6-(1,1...	340	C23H32O2	000119-47-1	87	
2	Phenol, 2,2'-methylenebis[6-(1,1...	340	C23H32O2	000119-47-1	86	
3	Phenol, 2,2'-methylenebis[6-(1,1...	340	C23H32O2	000119-47-1	70	
4	Benzofuran-2-one, 4-amino-2,3-di...	177	C10H11NO2	1000129-51-7	38	
5	2,3,5,6-Tetramethylbenzamide	177	C11H15NO	099858-56-7	38	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,2'-meth...	20.56	16.6	ng/ul	708868	5	21.16	853202	20.0