

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091923\
 Data File : BM041904.D
 Acq On : 19 Sep 2023 13:11
 Operator : MA/JU
 Sample : 04346-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WP

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM041904.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.498	385	393	404	rBV	7046785	10324966	48.25%	3.770%
2	7.116	658	668	670	rBV	6234652	12324170	57.59%	4.500%
3	7.140	670	672	686	rVB	6106518	7668427	35.83%	2.800%
4	7.516	728	736	747	rBV	6520671	10488621	49.01%	3.830%
5	7.957	804	811	822	rVB	1090895	1717937	8.03%	0.627%
6	8.398	877	886	903	rBV	7988295	13182635	61.60%	4.814%
7	8.722	932	941	956	rBV	7312151	12066060	56.39%	4.406%
8	9.157	1002	1015	1025	rBV	4373561	7285121	34.04%	2.660%
9	9.898	1132	1141	1143	rBV	2134354	3486347	16.29%	1.273%
10	9.933	1143	1147	1168	rVB	8205536	14044709	65.63%	5.128%
11	10.416	1220	1229	1243	rBV	8024360	13369458	62.48%	4.882%
12	10.780	1283	1291	1299	rBV	1510783	2500125	11.68%	0.913%
13	10.951	1306	1320	1330	rBV	3516624	5895442	27.55%	2.153%
14	12.051	1496	1507	1523	rBV	8422807	14160695	66.17%	5.171%
15	13.039	1667	1675	1680	rBV	9414257	14712826	68.75%	5.372%
16	13.104	1680	1686	1700	rVV	7791168	11586116	54.14%	4.231%
17	13.227	1700	1707	1716	rVB	9923909	15091004	70.52%	5.511%
18	13.910	1817	1823	1830	rBV	177987	248608	1.16%	0.091%
19	14.610	1935	1942	1951	rBV	2291387	3296774	15.41%	1.204%
20	14.757	1958	1967	1986	rBV	6918412	10395583	48.58%	3.796%
21	15.227	2039	2047	2064	rBV	11563480	17136603	80.08%	6.257%
22	15.745	2127	2135	2153	rBV	6925748	10238345	47.84%	3.739%
23	16.104	2189	2196	2212	rBV	8263516	12047902	56.30%	4.399%
24	16.992	2340	2347	2366	rBV	11401097	17087784	79.85%	6.240%
25	17.362	2404	2410	2417	rBV	2790027	3830192	17.90%	1.399%
26	19.974	2848	2854	2859	rBV	16128944	21399291	100.00%	7.814%
27	21.539	3115	3120	3126	rBV	3145423	4082250	19.08%	1.491%
28	23.962	3525	3532	3544	rVB	2101143	4190396	19.58%	1.530%

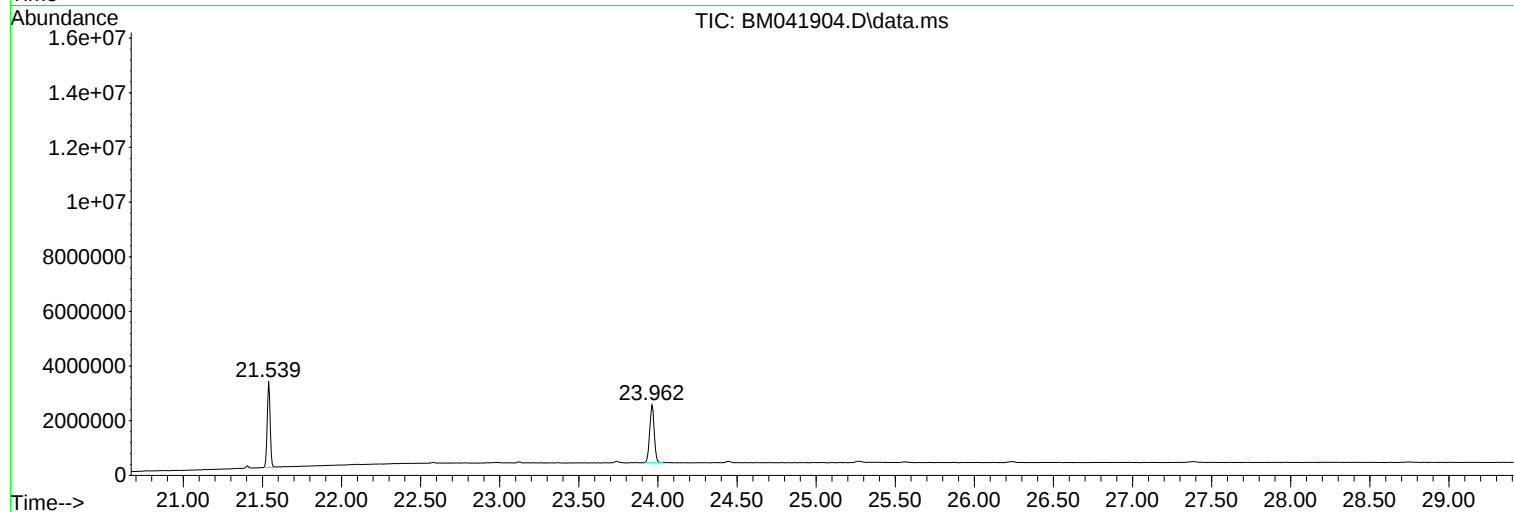
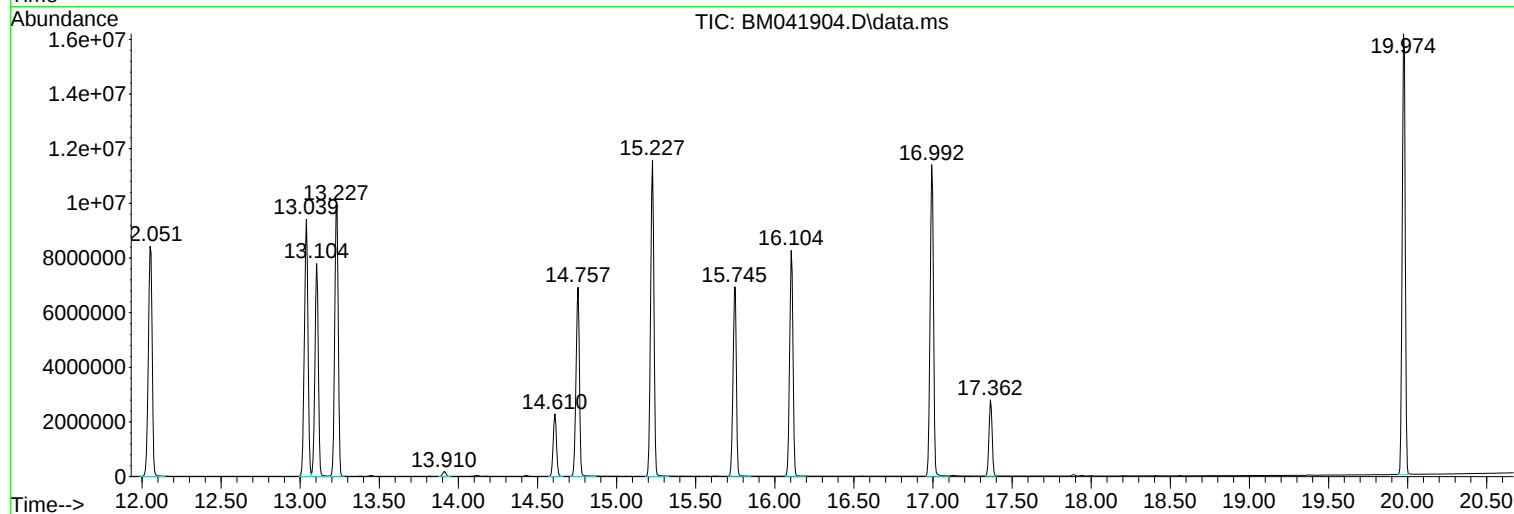
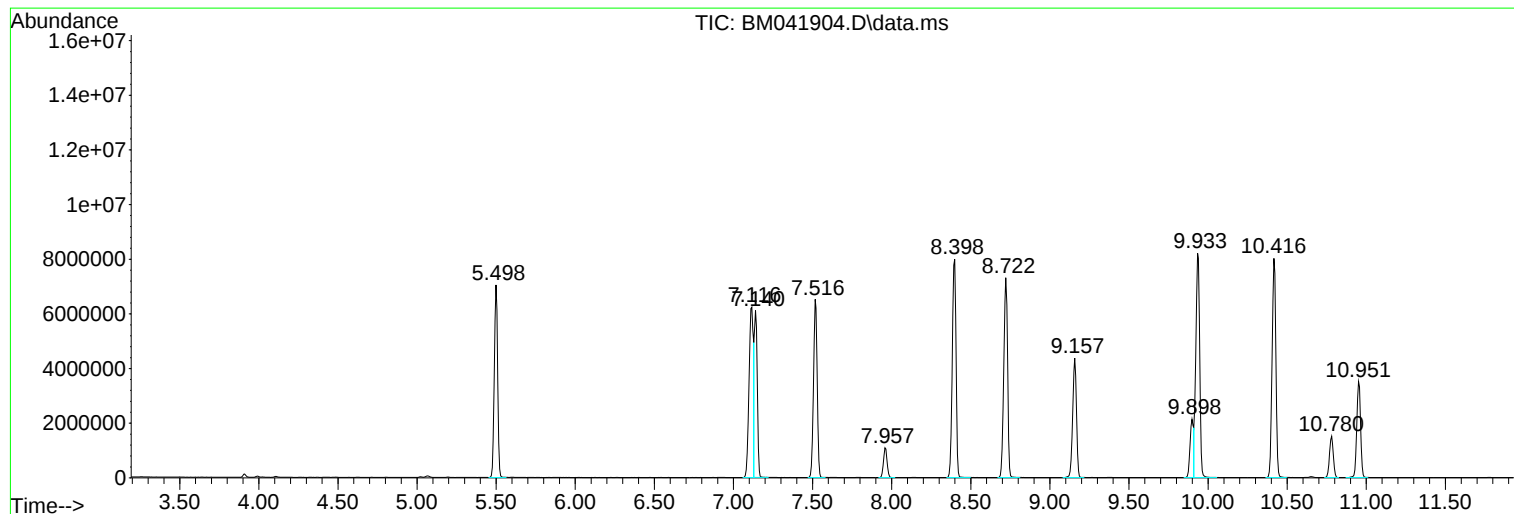
Sum of corrected areas: 273858387

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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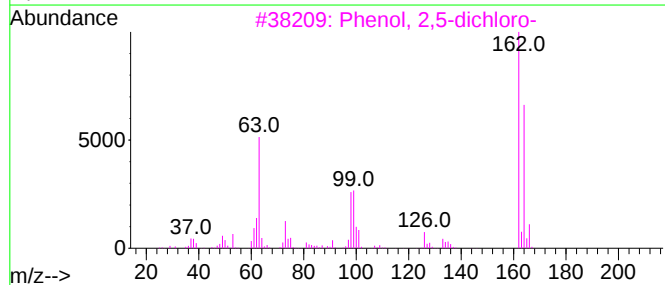
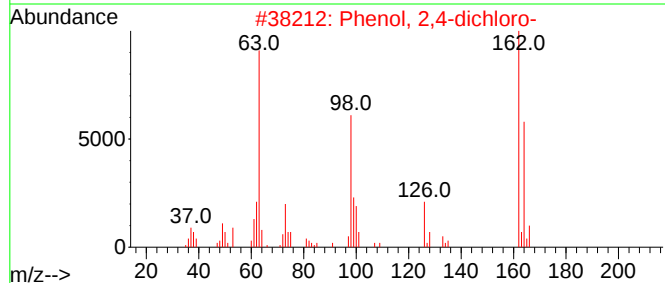
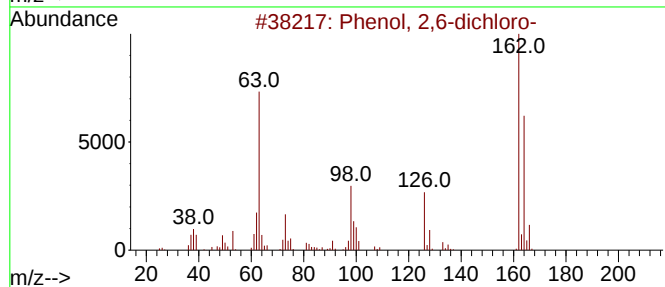
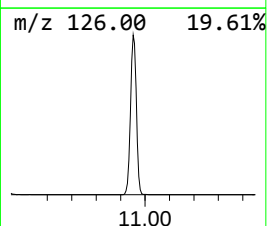
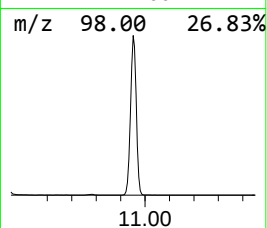
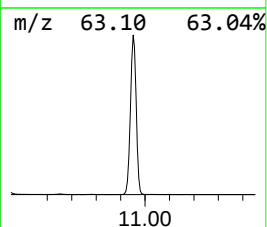
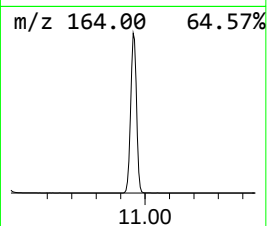
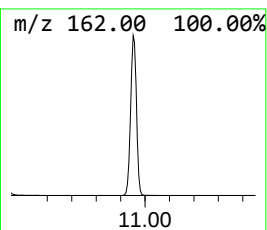
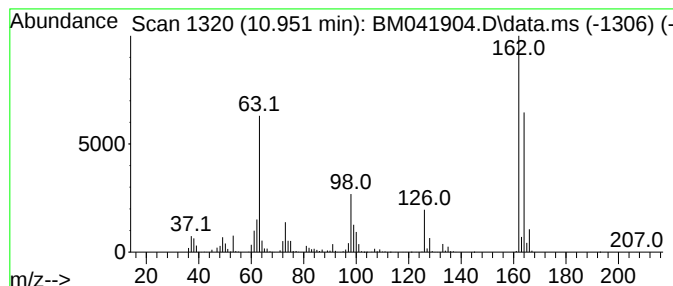
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenol, 2,6-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.951	47.16 ng	5895440	Naphthalene-d8	10.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	98
2			Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	97
3			Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	94
4			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	93
5			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-dic...	10.951	47.2	ng	5895440	2	10.781	2500130	20.0