

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
 Data File : BP021303.D
 Acq On : 02 Aug 2024 00:46
 Operator : CG/JU
 Sample : P3398-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E08K3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021303.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.163	26	30	40	rVB	36384	51297	1.08%	0.125%
2	3.593	100	103	122	rBV	129025	246398	5.20%	0.602%
3	3.881	147	152	167	rVB	763624	1066862	22.50%	2.607%
4	4.252	212	215	220	rVB7	3475	5424	0.11%	0.013%
5	5.916	493	498	506	rBV7	2974	7095	0.15%	0.017%
6	6.857	651	658	672	rBV	91996	279027	5.89%	0.682%
7	6.975	672	678	698	rVB	453125	784518	16.55%	1.917%
8	7.175	705	712	734	rBV	529547	1011163	21.33%	2.471%
9	7.622	781	788	805	rBV	591111	1071235	22.59%	2.618%
10	8.357	907	913	938	rBV2	345253	854520	18.02%	2.088%
11	8.793	977	987	1006	rBV	507978	1023210	21.58%	2.500%
12	8.928	1008	1010	1021	rVB9	4961	11625	0.25%	0.028%
13	9.504	1100	1108	1128	rBV	422687	900579	18.99%	2.201%
14	10.051	1194	1201	1232	rBV	498616	1314717	27.73%	3.213%
15	10.387	1251	1258	1271	rBV	885457	1443229	30.44%	3.527%
16	10.563	1279	1288	1316	rBV	392682	1141007	24.07%	2.788%
17	11.263	1402	1407	1408	rBV5	4123	5778	0.12%	0.014%
18	12.016	1529	1535	1547	rBV3	9245	28008	0.59%	0.068%
19	12.881	1678	1682	1686	rVV5	4253	6228	0.13%	0.015%
20	13.110	1713	1721	1732	rBV	118209	220847	4.66%	0.540%
21	13.239	1739	1743	1748	rVB8	2996	5306	0.11%	0.013%
22	13.681	1810	1818	1833	rBV	1082757	1794938	37.86%	4.386%
23	13.957	1858	1865	1885	rBV	1463421	2231366	47.06%	5.452%
24	14.263	1910	1917	1931	rBV	1264999	1893039	39.93%	4.626%
25	14.657	1976	1984	1988	rBV3	33721	78203	1.65%	0.191%
26	15.063	2050	2053	2058	rVB7	3354	4992	0.11%	0.012%
27	15.280	2083	2090	2106	rBV	1540482	2826267	59.61%	6.906%
28	15.463	2115	2121	2141	rBV	310546	796841	16.81%	1.947%
29	15.916	2194	2198	2202	rBV2	4017	5855	0.12%	0.014%
30	16.880	2357	2362	2366	rBV4	4235	7396	0.16%	0.018%
31	17.092	2391	2398	2407	rBV	1350660	2153603	45.42%	5.262%
32	17.198	2408	2416	2435	rVV2	1716182	3070101	64.75%	7.502%
33	18.027	2552	2557	2566	rBV	41038	82487	1.74%	0.202%
34	19.127	2740	2744	2749	rBV3	32155	50076	1.06%	0.122%
35	19.210	2754	2758	2764	rBV	29498	48665	1.03%	0.119%
36	19.368	2781	2785	2789	rBV6	28657	48091	1.01%	0.118%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

37	19.580	2814	2821	2841	rVB	2741174	3793455	80.01%	9.270%
38	21.545	3149	3155	3166	rBV2	1479444	2524561	53.25%	6.169%
39	24.603	3666	3675	3691	rVB	1977962	4741160	100.00%	11.585%
40	24.839	3706	3715	3730	rVB	1060351	3294662	69.49%	8.051%

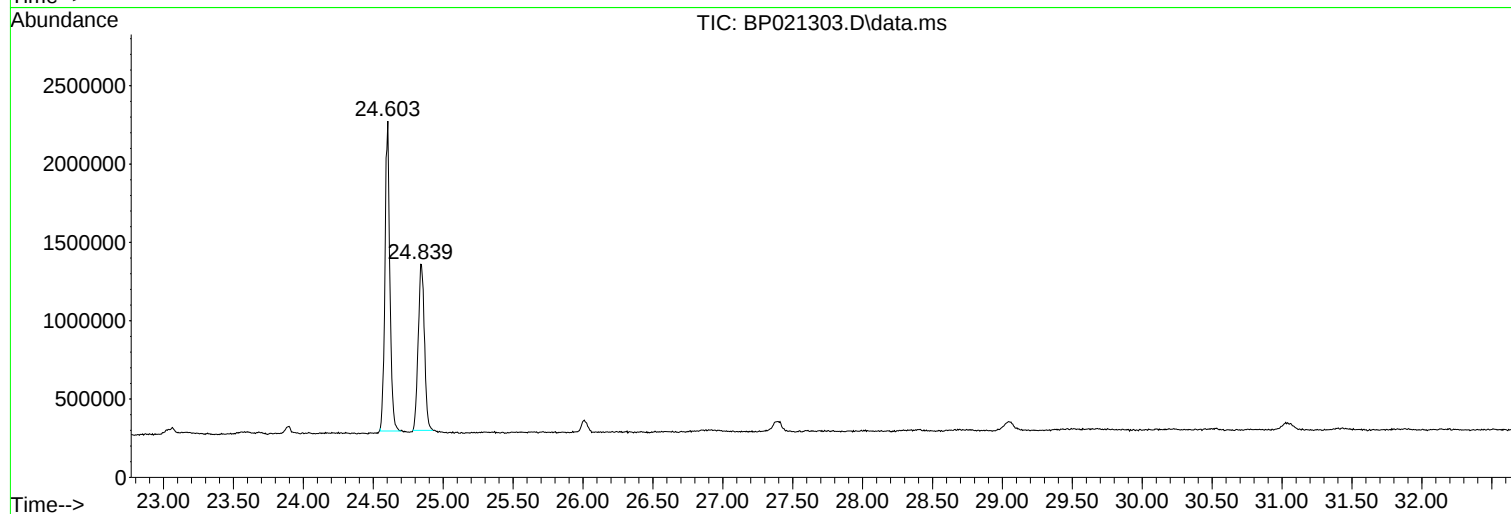
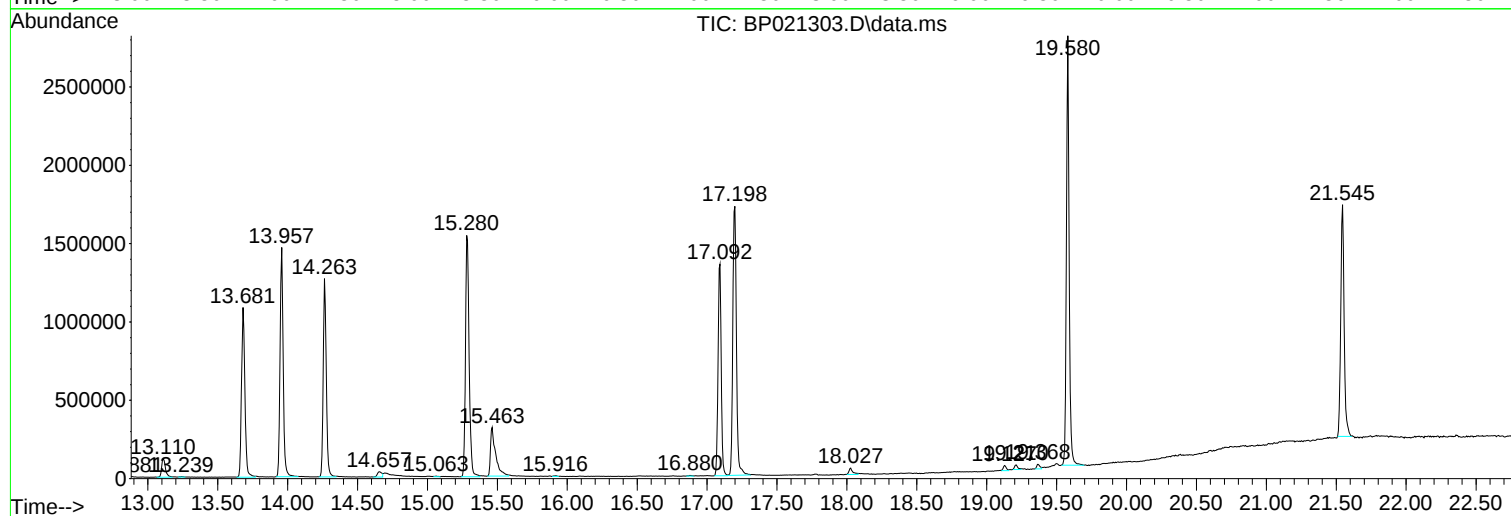
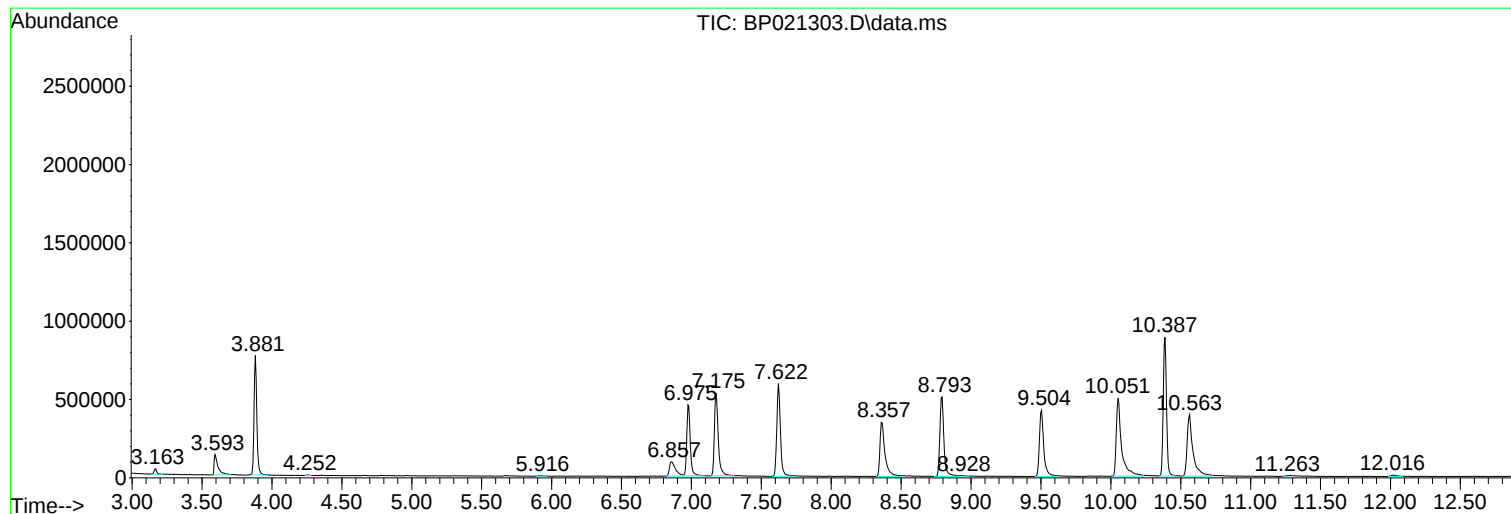
Sum of corrected areas: 40923831

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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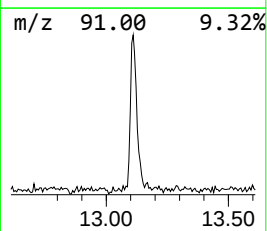
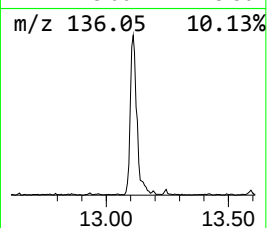
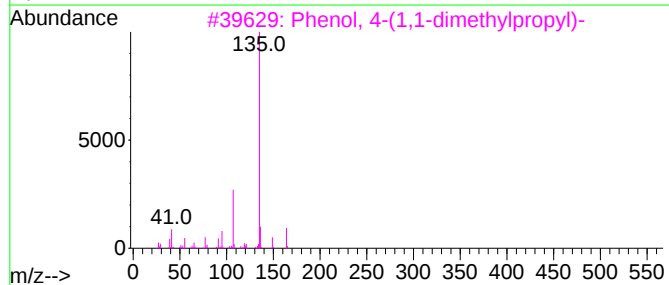
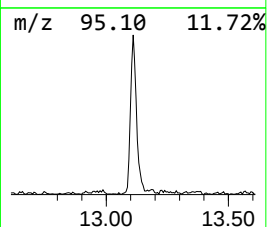
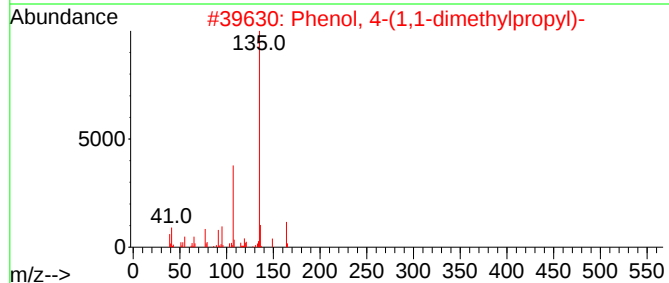
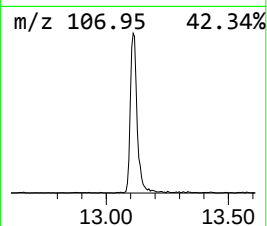
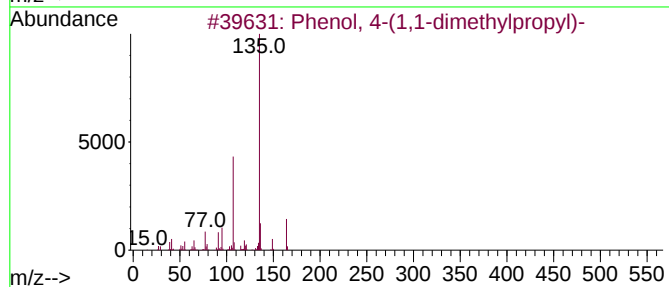
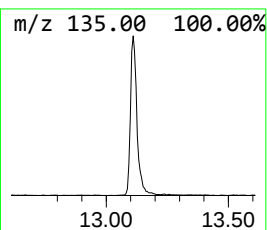
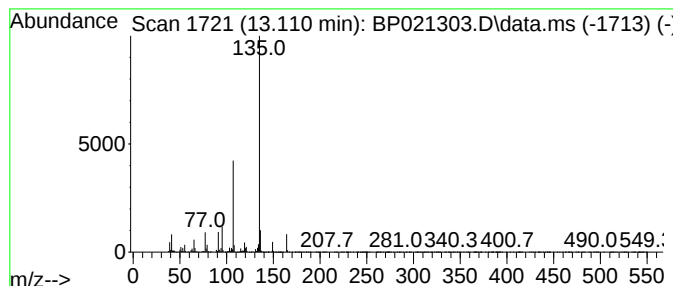
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 4-(1,1-dimethylprop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.110	2.33 ng/ul	220847	Acenaphthene-d10	14.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	95
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	91
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	87
4			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
5			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	81



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 4-(1,1-...	13.110	2.3	ng/ul	220847	3	14.263	1893040	20.0