

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019254.D
 Acq On : 03 Jan 2024 18:24
 Operator : MA/JU
 Sample : 05952-03DL 30X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCF88DL

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019254.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.681	99	101	111	rVB3	7670	12411	0.15%	0.058%
2	3.975	147	151	155	rBV2	10482	14406	0.17%	0.067%
3	6.993	658	664	673	rVB2	15267	31108	0.36%	0.144%
4	7.140	685	689	694	rVB2	14350	20830	0.24%	0.097%
5	7.334	717	722	728	rVB	17779	29339	0.34%	0.136%
6	7.799	792	801	809	rBV	657911	1058940	12.40%	4.917%
7	7.875	809	814	820	rVB3	6665	15317	0.18%	0.071%
8	8.534	921	926	932	rBV3	14817	31272	0.37%	0.145%
9	8.969	992	1000	1008	rBV2	12012	27754	0.33%	0.129%
10	9.687	1117	1122	1129	rVB6	9696	14947	0.18%	0.069%
11	10.246	1211	1217	1225	rBV3	11427	29895	0.35%	0.139%
12	10.593	1264	1276	1286	rBV	792685	1318744	15.44%	6.124%
13	12.028	1509	1520	1521	rBV	6357	16391	0.19%	0.076%
14	12.681	1627	1631	1639	rVB10	9833	16476	0.19%	0.077%
15	13.345	1739	1744	1748	rVB8	6991	12181	0.14%	0.057%
16	13.704	1800	1805	1809	rBV8	5795	10523	0.12%	0.049%
17	13.863	1827	1832	1838	rVB	31129	46815	0.55%	0.217%
18	14.134	1873	1878	1886	rVB	31850	53703	0.63%	0.249%
19	14.269	1897	1901	1905	rVB7	5792	9953	0.12%	0.046%
20	14.339	1908	1913	1919	rVB3	15376	24582	0.29%	0.114%
21	14.439	1924	1930	1937	rBV	1079120	1550034	18.15%	7.198%
22	14.669	1965	1969	1974	rVB	109569	143866	1.68%	0.668%
23	14.992	2022	2024	2028	rVB4	9556	12015	0.14%	0.056%
24	15.051	2028	2034	2035	rBV4	35885	51779	0.61%	0.240%
25	15.075	2035	2038	2046	rVB	82967	131620	1.54%	0.611%
26	15.275	2068	2072	2078	rBV4	30166	53805	0.63%	0.250%
27	15.439	2096	2100	2104	rVB	47788	59840	0.70%	0.278%
28	15.498	2106	2110	2113	rBV6	18836	36457	0.43%	0.169%
29	16.851	2334	2340	2347	rBV	5742233	8538588	100.00%	39.650%
30	16.998	2360	2365	2375	rVB2	598577	1053114	12.33%	4.890%
31	17.198	2394	2399	2404	rBV	1336432	1865657	21.85%	8.663%
32	21.304	3092	3097	3107	rVB	1932292	2430208	28.46%	11.285%
33	23.668	3493	3499	3509	rVB	1376737	2812336	32.94%	13.059%

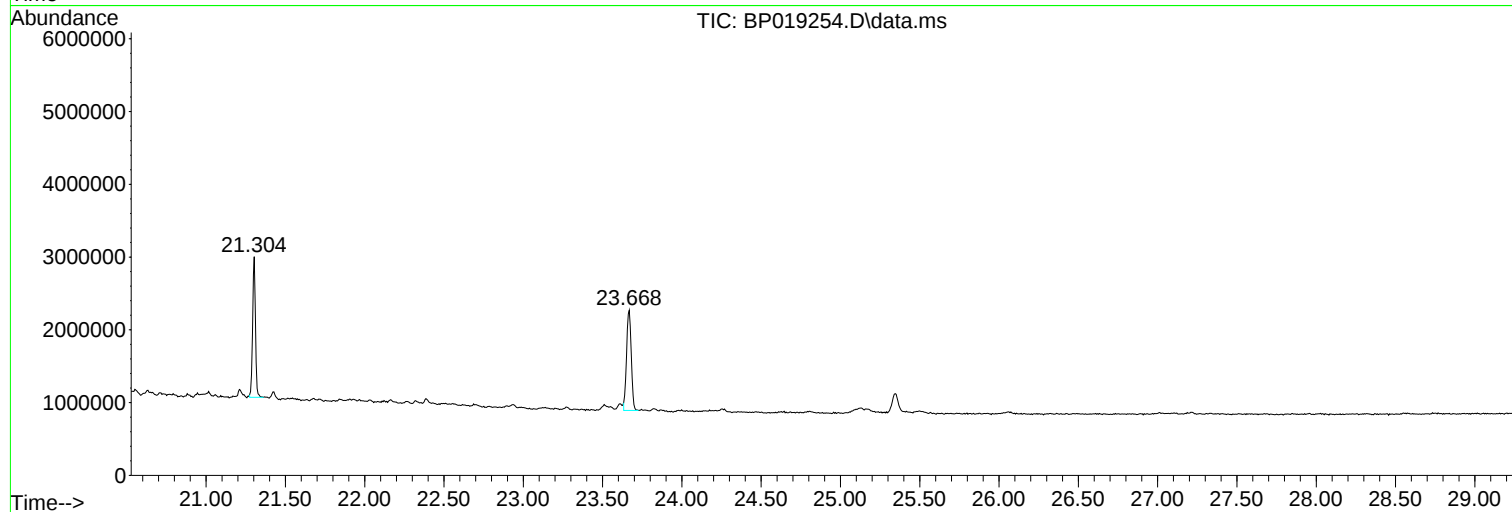
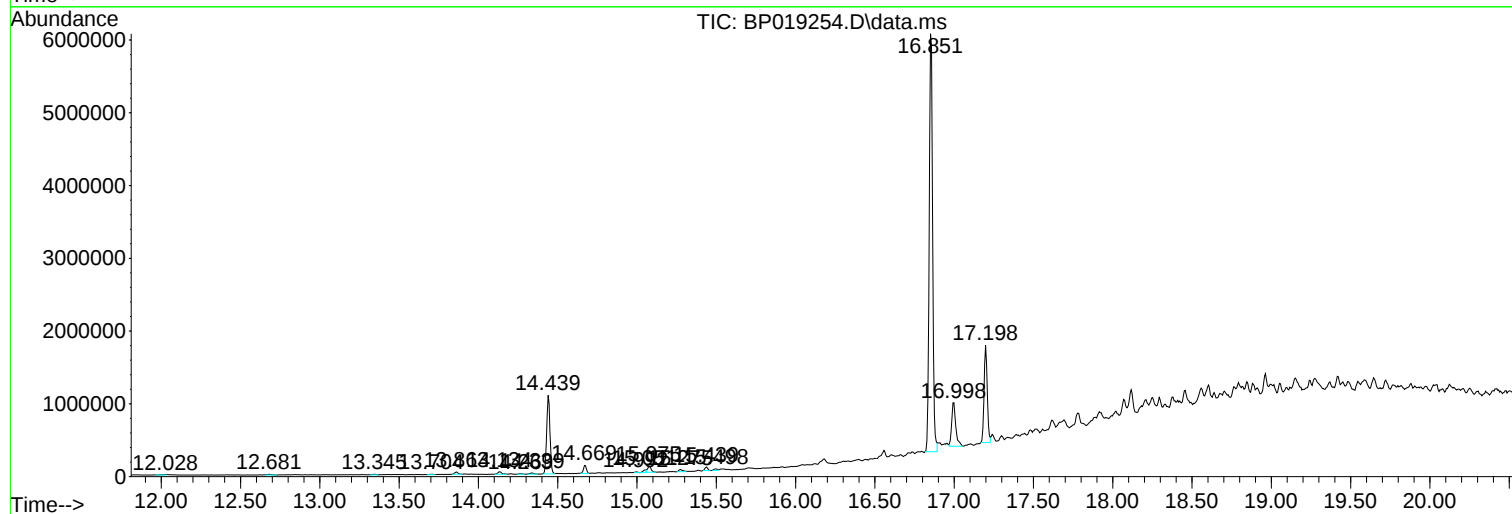
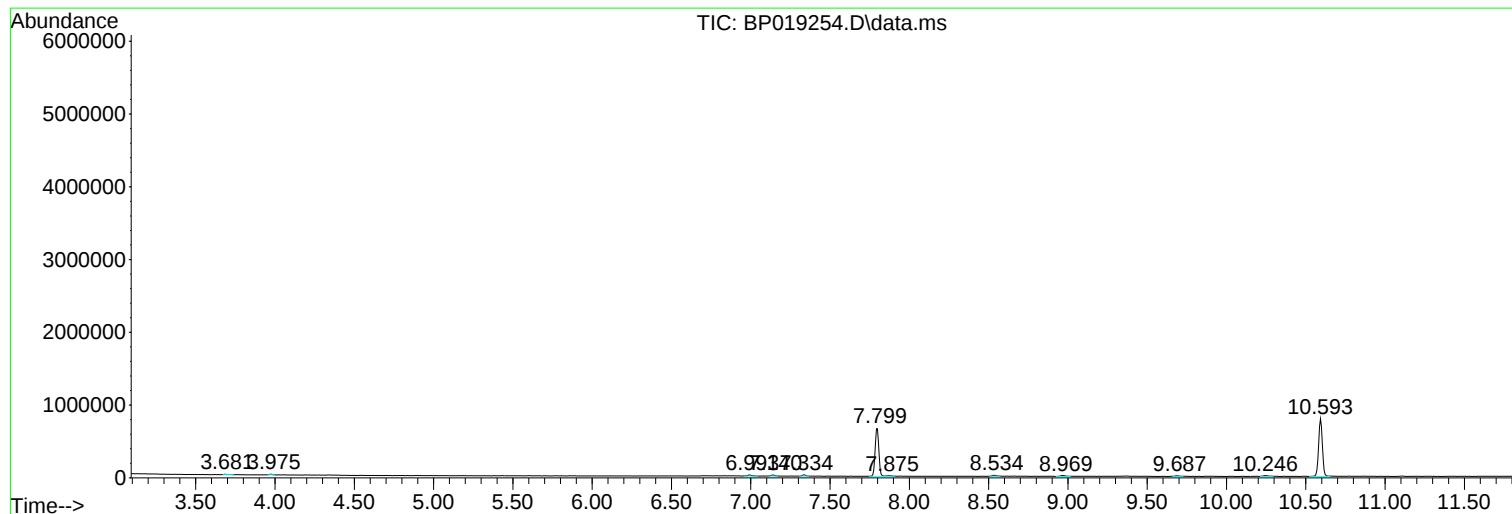
Sum of corrected areas: 21534906

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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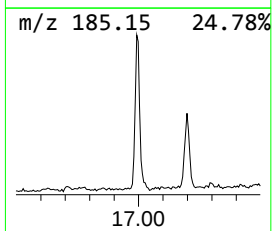
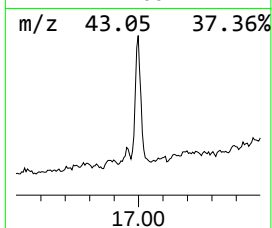
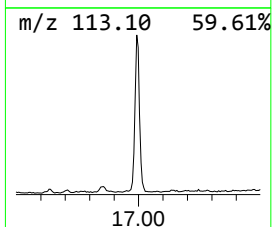
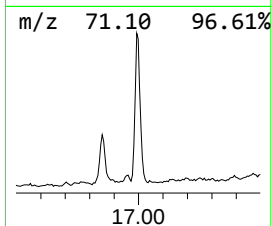
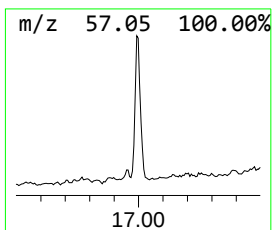
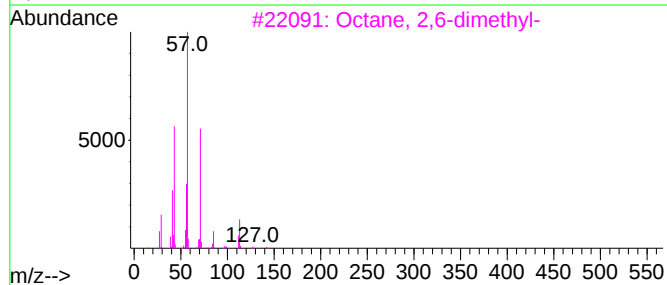
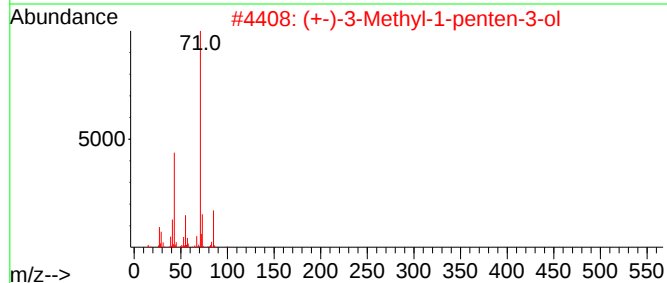
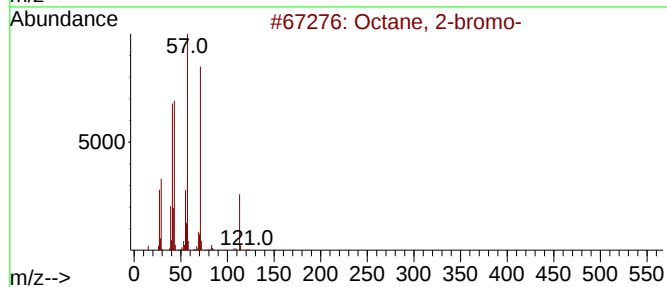
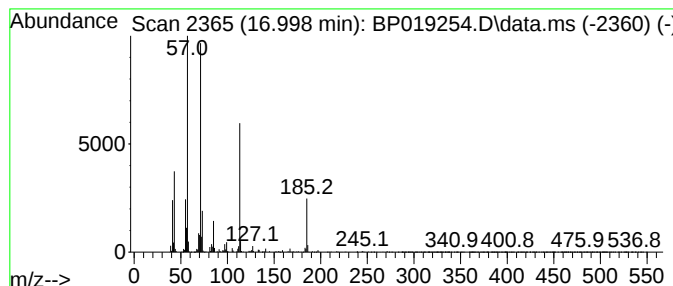
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.998	11.29 ng/ul	1053110	Phenanthrene-d10	17.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-bromo-			192	C8H17Br	000557-35-7	47
2	(+)-3-Methyl-1-penten-3-ol			100	C6H12O	1010238-75-4	46
3	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	43
4	Dodecane, 4,6-dimethyl-			198	C14H30	061141-72-8	43
5	Sulfurous acid, 2-ethylhexyl iso...			250	C12H26O3S	1000309-18-7	38



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	16.998	11.3	ng/ul	1053110	4	17.198	1865660	20.0