

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
 Data File : BP021314.D
 Acq On : 02 Aug 2024 09:33
 Operator : CG/JU
 Sample : P3397-08
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHCM8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021314.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	27	30	39	rVB	14051	23171	0.72%	0.096%
2	3.634	106	110	120	rBV3	9754	23096	0.72%	0.096%
3	3.881	148	152	158	rVB	14724	20083	0.63%	0.083%
4	6.858	651	658	673	rBV	136094	380325	11.86%	1.581%
5	6.981	673	679	696	rVB	161541	305192	9.52%	1.269%
6	7.181	705	713	730	rBV	219652	428802	13.37%	1.783%
7	7.622	780	788	805	rBV	504389	866426	27.02%	3.602%
8	8.369	908	915	939	rBV	219791	510762	15.93%	2.124%
9	8.793	979	987	1007	rBV	198098	394388	12.30%	1.640%
10	9.363	1078	1084	1089	rBV9	3287	6905	0.22%	0.029%
11	9.504	1101	1108	1126	rBV	148519	332761	10.38%	1.384%
12	10.069	1194	1204	1227	rBV	141252	499296	15.57%	2.076%
13	10.387	1250	1258	1275	rBV	562018	1168452	36.44%	4.858%
14	10.563	1280	1288	1304	rBV	156909	452997	14.13%	1.883%
15	10.787	1321	1326	1330	rVB7	2080	4266	0.13%	0.018%
16	11.193	1389	1395	1396	rBV	8739	11173	0.35%	0.046%
17	11.210	1396	1398	1407	rVB7	7431	17416	0.54%	0.072%
18	12.010	1527	1534	1539	rBV7	4677	9064	0.28%	0.038%
19	13.675	1811	1817	1837	rBV	450450	722951	22.54%	3.006%
20	13.957	1857	1865	1879	rVV	591965	960124	29.94%	3.992%
21	14.051	1879	1881	1887	rVB6	2955	4661	0.15%	0.019%
22	14.263	1910	1917	1935	rBV	963791	1552193	48.40%	6.454%
23	14.492	1946	1956	1961	rBV4	16731	43893	1.37%	0.182%
24	14.645	1975	1982	1983	rBV	45042	69436	2.17%	0.289%
25	15.122	2060	2063	2067	rVB6	3327	4230	0.13%	0.018%
26	15.287	2084	2091	2106	rBV	767940	1249727	38.97%	5.196%
27	15.416	2111	2113	2118	rVB5	2793	3223	0.10%	0.013%
28	15.481	2118	2124	2144	rBV2	93561	232170	7.24%	0.965%
29	16.828	2350	2353	2356	rBV5	2509	3686	0.11%	0.015%
30	16.869	2358	2360	2365	rVB6	4645	6579	0.21%	0.027%
31	17.086	2389	2397	2409	rBV	1195205	1946956	60.71%	8.095%
32	17.192	2409	2415	2430	rVV	899424	1388802	43.31%	5.774%
33	17.369	2442	2445	2450	rVB7	3728	6187	0.19%	0.026%
34	17.757	2509	2511	2519	rVB9	6359	10120	0.32%	0.042%
35	18.016	2549	2555	2565	rBV4	40689	106193	3.31%	0.442%
36	19.198	2753	2756	2759	rBV3	14612	22981	0.72%	0.096%

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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

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Title : SVOA CALIBRATION

37	19.345	2778	2781	2786	rBV5	20687	32795	1.02%	0.136%
38	19.569	2813	2819	2835	rBV	1232539	1963892	61.24%	8.165%
39	21.169	3087	3091	3100	rBV2	138303	257755	8.04%	1.072%
40	21.527	3146	3152	3165	rBV2	1401515	2556628	79.72%	10.630%
41	24.598	3665	3674	3692	rVB	752793	2245060	70.01%	9.334%
42	24.821	3703	3712	3729	rVB	1070330	3206880	100.00%	13.333%

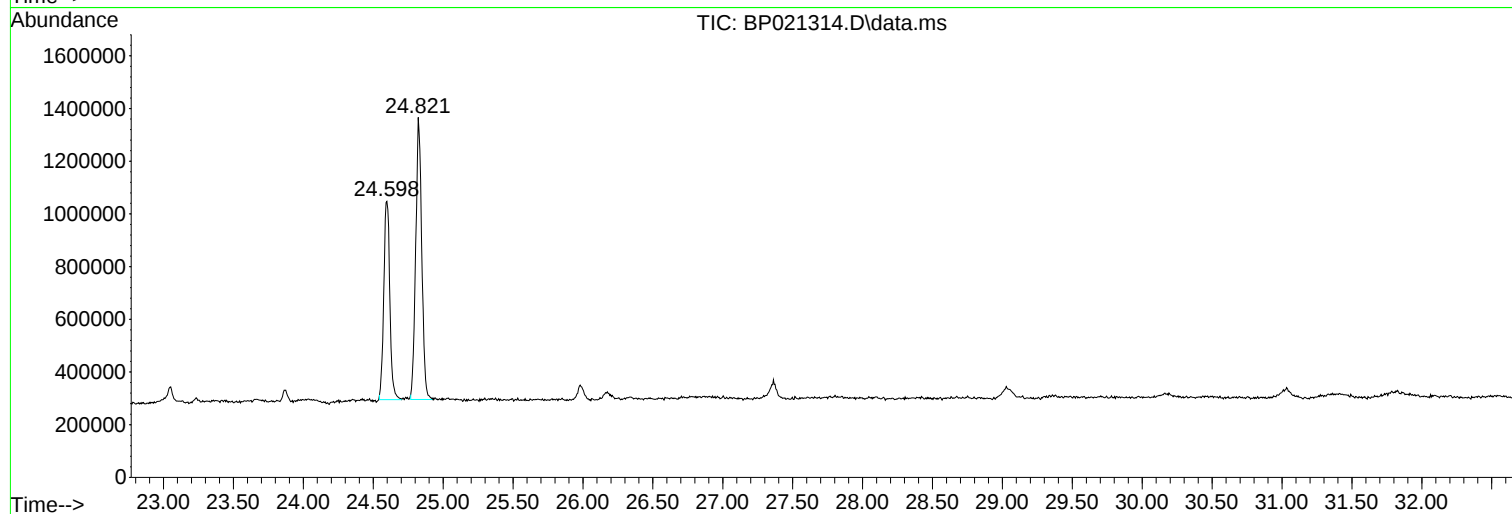
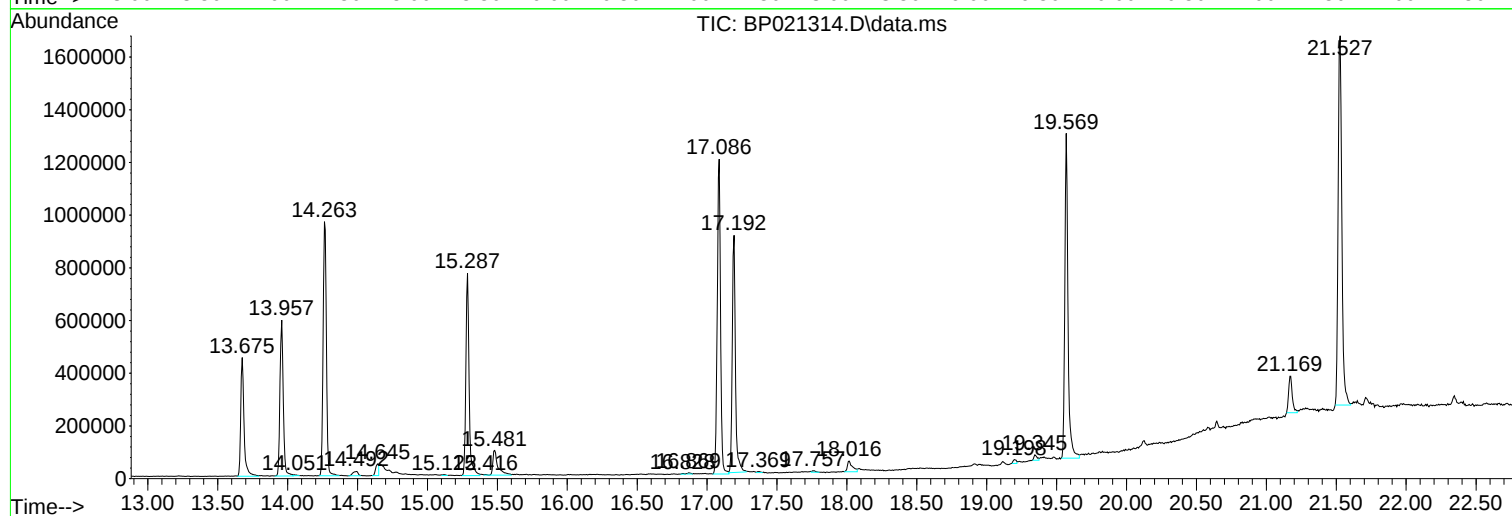
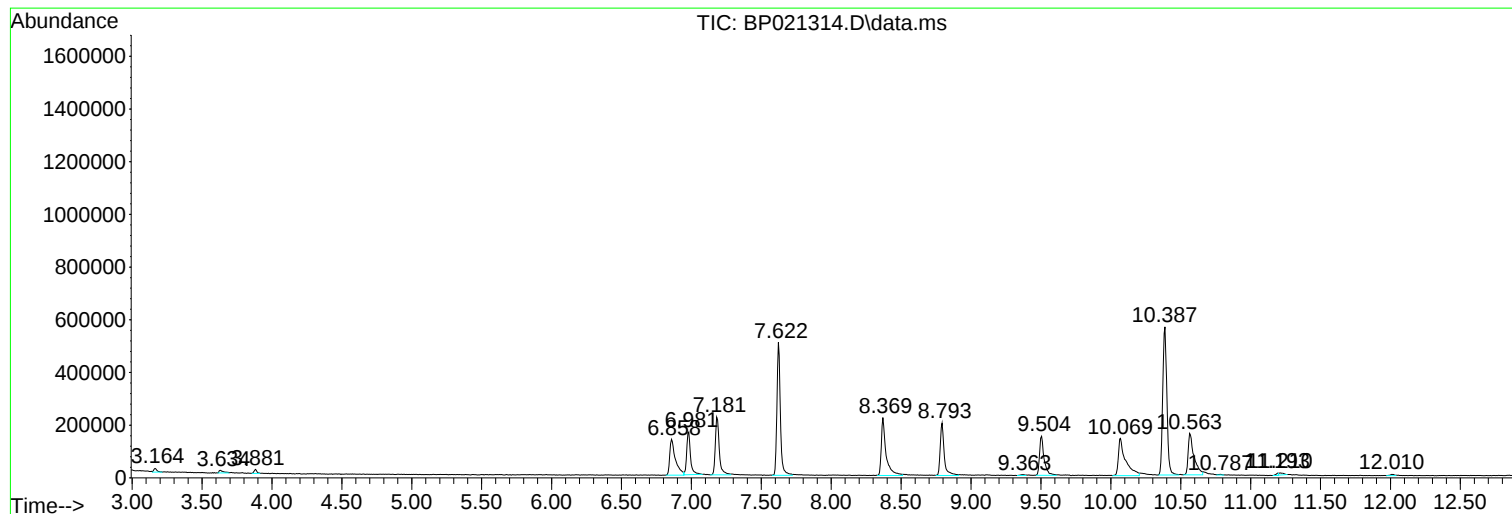
Sum of corrected areas: 24051697

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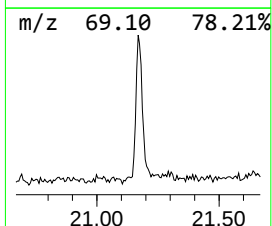
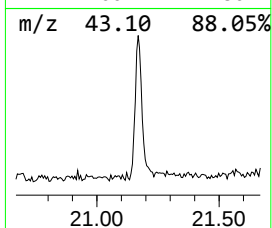
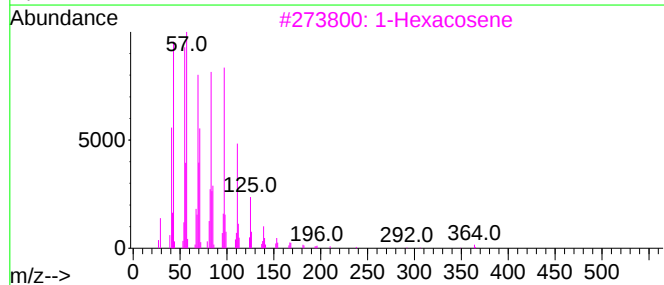
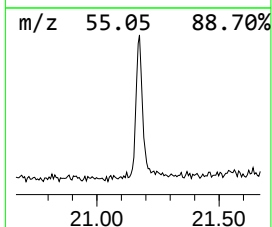
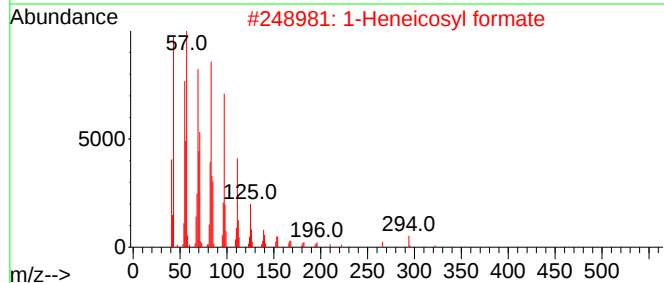
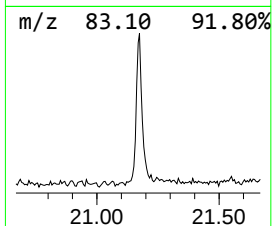
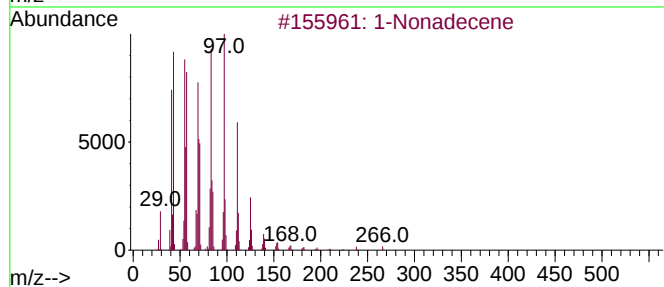
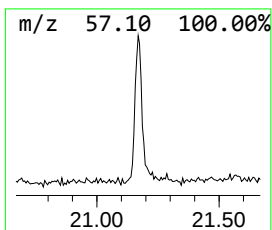
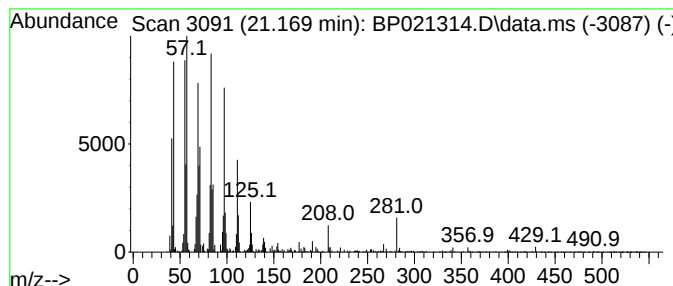
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.169	2.02 ng/ul	257755	Chrysene-d12	21.527

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene		266	C19H38	018435-45-5	98
2	1-Heneicosyl formate		340	C22H44O2	077899-03-7	97
3	1-Hexacosene		364	C26H52	018835-33-1	94
4	n-Nonadecanol-1		284	C19H40O	001454-84-8	94
5	Pentafluoropropionic acid, penta...		374	C18H31F5O2	959092-08-1	94



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	21.169	2.0	ng/u1	257755	5	21.527	2556630	20.0