

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
 Data File : BP020964.D
 Acq On : 15 Jul 2024 18:53
 Operator : MA/JU
 Sample : P3087-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CB7

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: BP020964.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	12	15	23	rVB	127796	189728	3.42%	0.249%
2	3.252	41	45	53	rVB	38107	51169	0.92%	0.067%
3	3.540	91	94	99	rBV	15243	22371	0.40%	0.029%
4	3.687	118	119	127	rVV	16275	28383	0.51%	0.037%
5	3.764	127	132	140	rVB	73445	105976	1.91%	0.139%
6	3.969	163	167	172	rVB	17982	26768	0.48%	0.035%
7	4.522	256	261	267	rVB4	19875	33647	0.61%	0.044%
8	4.869	314	320	328	rVB2	39676	77696	1.40%	0.102%
9	4.975	334	338	345	rVB3	11108	18850	0.34%	0.025%
10	6.810	637	650	654	rBV8	6271	18967	0.34%	0.025%
11	6.899	658	665	678	rBV	602520	1099243	19.82%	1.444%
12	7.046	683	690	701	rVB	508920	865632	15.61%	1.137%
13	7.246	717	724	732	rBV	764933	1174581	21.18%	1.543%
14	7.693	790	800	808	rBV	643632	1079132	19.46%	1.418%
15	8.410	915	922	935	rBV	678553	1229981	22.18%	1.616%
16	8.516	935	940	950	rVB2	25512	55635	1.00%	0.073%
17	8.846	988	996	1011	rBV	645787	1148273	20.70%	1.509%
18	8.987	1016	1020	1027	rVB9	7595	15994	0.29%	0.021%
19	9.551	1103	1116	1130	rBV	571974	1008261	18.18%	1.325%
20	10.098	1201	1209	1225	rBV	758519	1564918	28.22%	2.056%
21	10.457	1263	1270	1276	rBV	876212	1508952	27.21%	1.982%
22	10.504	1276	1278	1284	rVB	42094	60071	1.08%	0.079%
23	10.604	1288	1295	1308	rBV	309909	690082	12.44%	0.907%
24	11.204	1391	1397	1405	rBV2	19995	59785	1.08%	0.079%
25	12.045	1534	1540	1546	rVB3	13583	24774	0.45%	0.033%
26	12.110	1546	1551	1557	rBV	17343	29397	0.53%	0.039%
27	12.340	1585	1590	1596	rVB3	12336	24144	0.44%	0.032%
28	13.492	1777	1786	1791	rBV7	15459	39929	0.72%	0.052%
29	13.622	1802	1808	1813	rBV2	19600	37352	0.67%	0.049%
30	13.710	1813	1823	1830	rBV	1577076	2320874	41.85%	3.049%
31	14.004	1866	1873	1882	rVV	1920135	2803520	50.55%	3.683%
32	14.192	1901	1905	1909	rVB3	12049	16304	0.29%	0.021%
33	14.310	1914	1925	1931	rBV	1410470	2098424	37.84%	2.757%
34	14.375	1931	1936	1943	rVV	130049	192503	3.47%	0.253%
35	14.439	1943	1947	1955	rVV2	33344	53857	0.97%	0.071%
36	14.534	1955	1963	1964	rVV3	41457	82880	1.49%	0.109%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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

37	14.569	1964	1969	1985	rVV	333245	890497	16.06%	1.170%
38	14.716	1989	1994	2005	rVV	114210	267479	4.82%	0.351%
39	14.798	2006	2008	2021	rVB6	21306	43548	0.79%	0.057%
40	15.116	2055	2062	2066	rBV	37001	63834	1.15%	0.084%
41	15.157	2066	2069	2073	rVB3	28516	35765	0.64%	0.047%
42	15.263	2082	2087	2090	rBV7	10322	20389	0.37%	0.027%
43	15.316	2090	2096	2102	rVV	2455560	3543350	63.89%	4.655%
44	15.375	2102	2106	2115	rVB2	134461	216873	3.91%	0.285%
45	15.463	2115	2121	2132	rBV	609712	1109388	20.00%	1.458%
46	15.575	2136	2140	2146	rVB3	91476	149713	2.70%	0.197%
47	15.675	2152	2157	2165	rVB	31357	62584	1.13%	0.082%
48	15.833	2180	2184	2191	rBV	35887	57798	1.04%	0.076%
49	16.045	2216	2220	2222	rBV	39786	57734	1.04%	0.076%
50	16.328	2263	2268	2274	rBV2	70602	105201	1.90%	0.138%
51	16.386	2275	2278	2280	rVV4	18661	25558	0.46%	0.034%
52	16.410	2280	2282	2287	rVB2	30409	35951	0.65%	0.047%
53	16.457	2287	2290	2295	rVB3	21455	28448	0.51%	0.037%
54	16.551	2304	2306	2312	rVB7	16402	26982	0.49%	0.035%
55	16.639	2316	2321	2325	rBV4	55329	86356	1.56%	0.113%
56	16.775	2339	2344	2350	rVB	72951	110156	1.99%	0.145%
57	16.839	2351	2355	2356	rBV2	23321	28372	0.51%	0.037%
58	16.863	2356	2359	2364	rVV4	52007	76135	1.37%	0.100%
59	16.933	2367	2371	2376	rVB	100248	143272	2.58%	0.188%
60	17.033	2384	2388	2392	rVB5	48081	70540	1.27%	0.093%
61	17.122	2396	2403	2407	rBV	1548768	2585244	46.61%	3.397%
62	17.169	2407	2411	2416	rVV	1814714	2627236	47.37%	3.452%
63	17.227	2416	2421	2425	rVV	2529247	3666497	66.11%	4.817%
64	17.257	2425	2426	2431	rVB	335673	283174	5.11%	0.372%
65	17.304	2431	2434	2439	rBV2	63326	92781	1.67%	0.122%
66	17.551	2467	2476	2480	rBV2	127512	284810	5.14%	0.374%
67	17.733	2501	2507	2514	rVB3	84200	182660	3.29%	0.240%
68	17.851	2524	2527	2534	rVB5	44718	79553	1.43%	0.105%
69	18.022	2552	2556	2558	rBV2	207676	277071	5.00%	0.364%
70	18.051	2558	2561	2564	rVV2	346034	436646	7.87%	0.574%
71	18.157	2576	2579	2584	rVB	84983	113185	2.04%	0.149%
72	18.222	2584	2590	2595	rBV3	374730	685144	12.35%	0.900%
73	18.263	2595	2597	2604	rVB2	135765	236642	4.27%	0.311%
74	18.333	2604	2609	2611	rBV2	61761	104346	1.88%	0.137%
75	18.545	2641	2645	2650	rVV	152737	215084	3.88%	0.283%
76	18.598	2650	2654	2661	rVB	201495	307719	5.55%	0.404%

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Peak separation: 5

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 Title : SVOA CALIBRATION

77	18.798	2685	2688	2693	rVB2	41515	61161	1.10%	0.080%
78	18.980	2715	2719	2722	rBV	82432	122093	2.20%	0.160%
79	19.251	2759	2765	2772	rBV	3334152	4811239	86.75%	6.321%
80	19.386	2784	2788	2796	rVB3	208288	383176	6.91%	0.503%
81	19.598	2819	2824	2827	rBV	3307583	4717528	85.06%	6.198%
82	19.633	2827	2830	2838	rVV	2243828	3073452	55.42%	4.038%
83	19.710	2839	2843	2849	rVB2	154488	212366	3.83%	0.279%
84	19.816	2858	2861	2866	rVB	155288	198775	3.58%	0.261%
85	20.145	2914	2917	2928	rVB2	435319	669603	12.07%	0.880%
86	20.251	2931	2935	2939	rBV2	266399	362448	6.54%	0.476%
87	20.310	2942	2945	2950	rVV2	179905	261447	4.71%	0.343%
88	20.463	2968	2971	2974	rBV	121534	140024	2.52%	0.184%
89	20.504	2976	2978	2984	rVB	147995	199673	3.60%	0.262%
90	20.939	3050	3052	3058	rVB	202113	285033	5.14%	0.374%
91	21.127	3081	3084	3088	rBV3	224085	283124	5.10%	0.372%
92	21.168	3089	3091	3095	rVB	209339	233473	4.21%	0.307%
93	21.216	3095	3099	3106	rBV3	651594	1087601	19.61%	1.429%
94	21.457	3135	3140	3145	rVB	2763043	3422779	61.71%	4.497%
95	21.557	3150	3157	3162	rVV3	2298747	5546198	100.00%	7.287%
96	21.604	3162	3165	3178	rVB	1482608	2307413	41.60%	3.032%
97	23.810	3534	3540	3548	rBV2	903513	2848839	51.37%	3.743%
98	23.945	3559	3563	3570	rVB	358006	652980	11.77%	0.858%
99	24.627	3672	3679	3685	rBV2	926644	2307081	41.60%	3.031%
100	24.868	3711	3720	3729	rBV3	900595	2634467	47.50%	3.461%

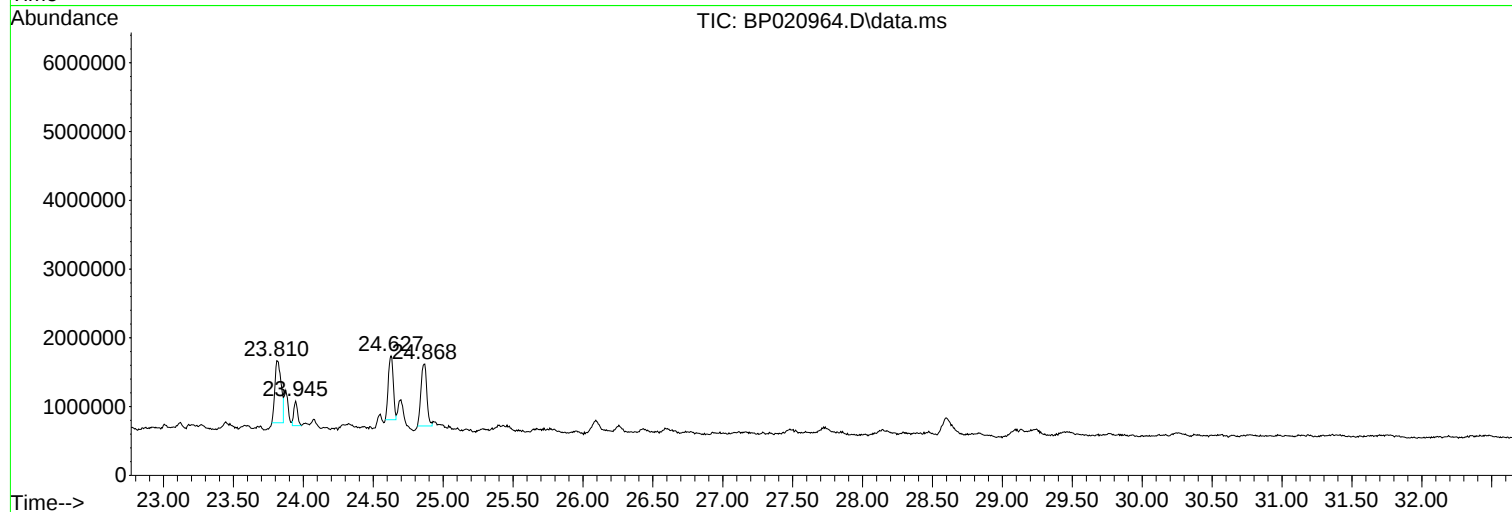
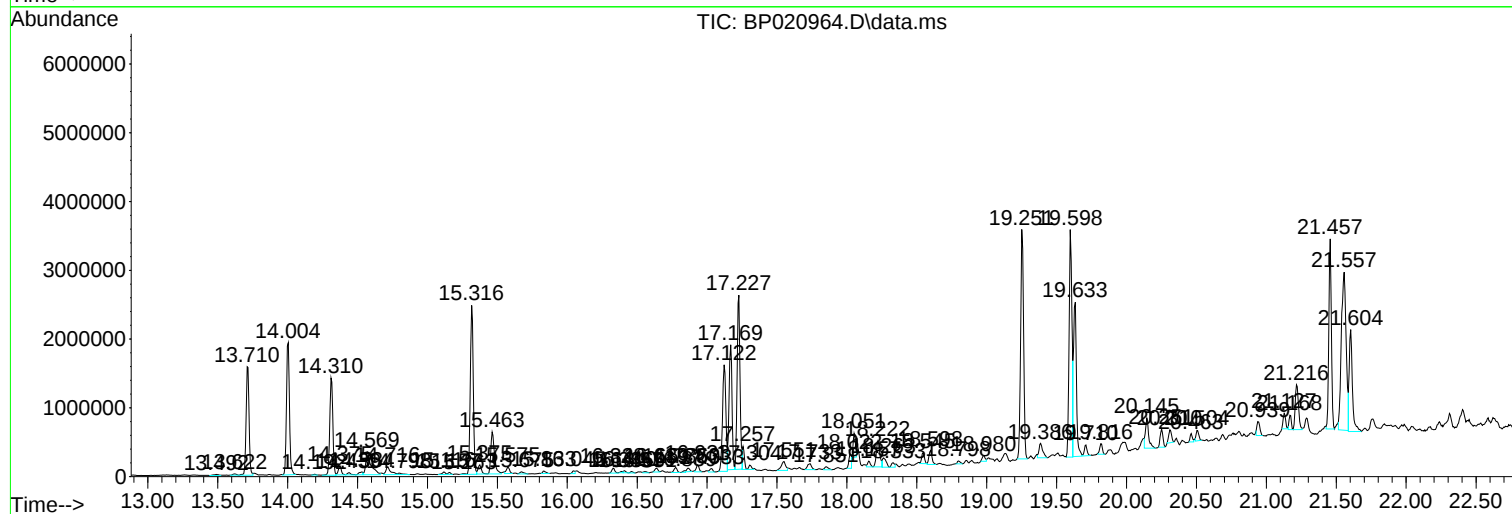
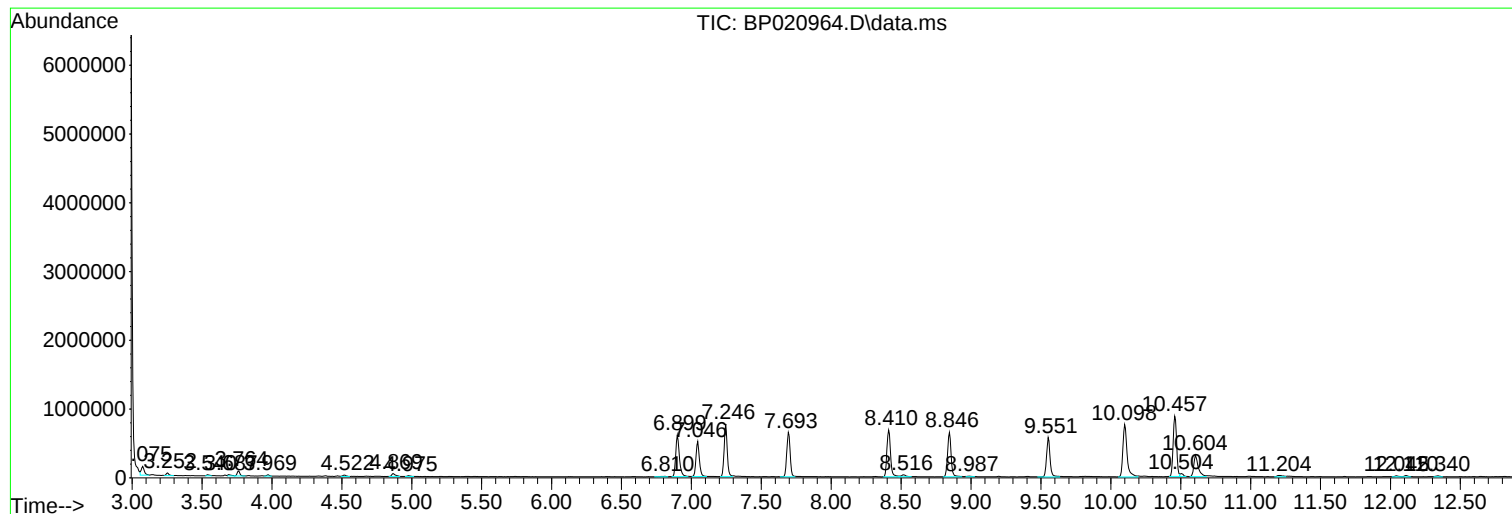
Sum of corrected areas: 76113741

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
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Instrument :
BNA_P
ClientSampleId :
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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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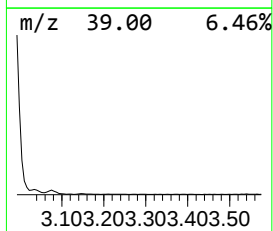
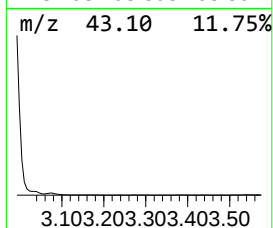
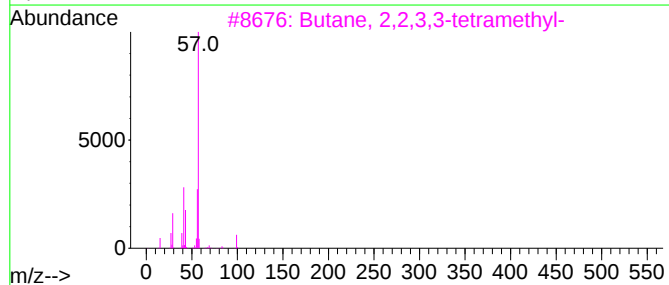
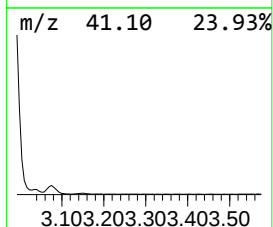
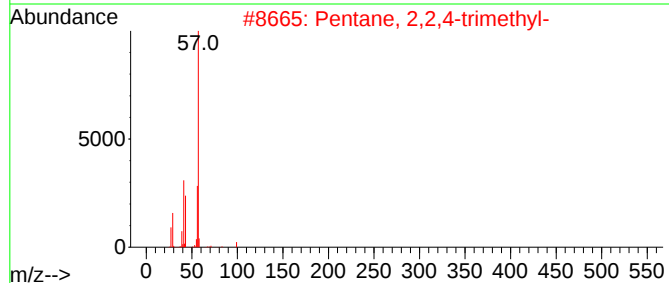
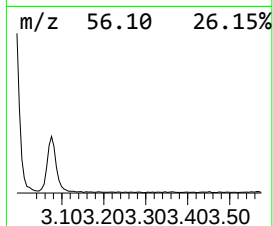
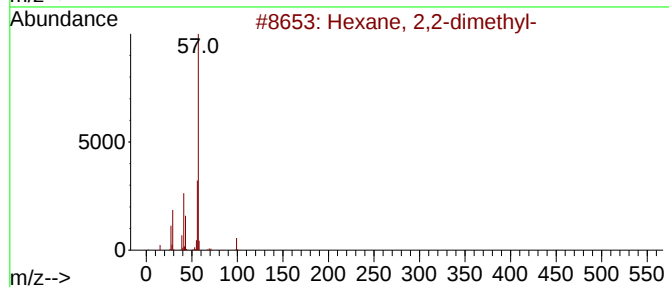
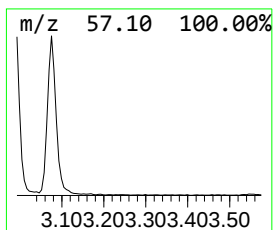
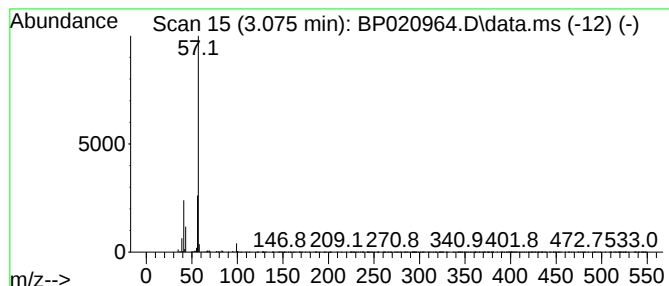
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	3.52 ng/ul	189728	1,4-Dichlorobenzene-d4	7.693

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
2		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	45
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	42
4		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	40
5		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	39



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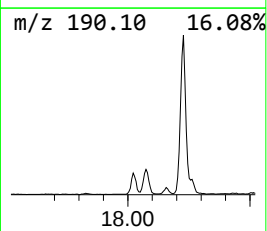
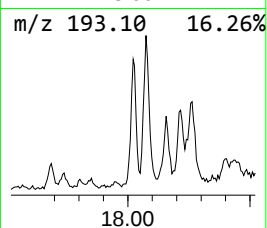
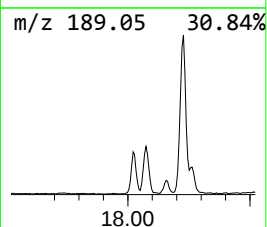
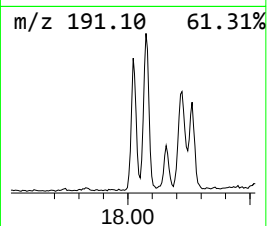
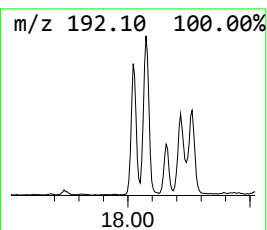
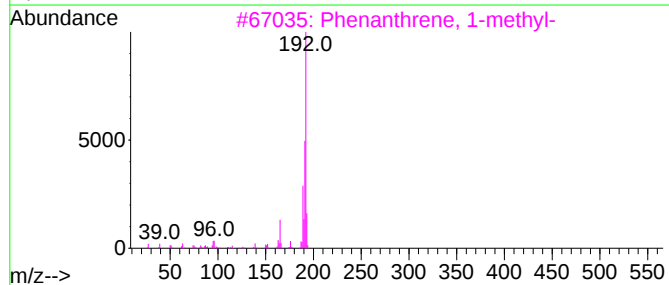
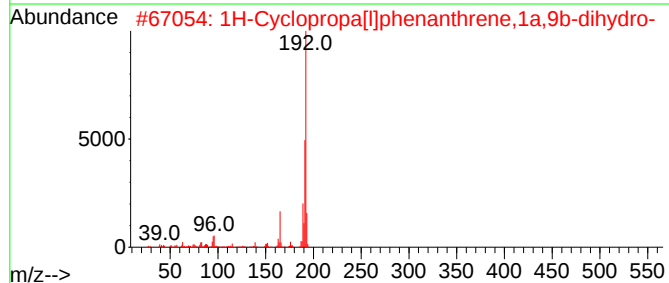
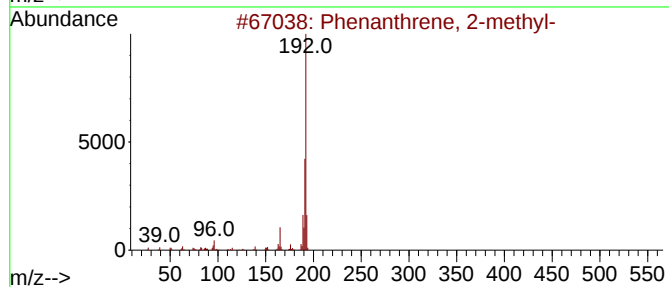
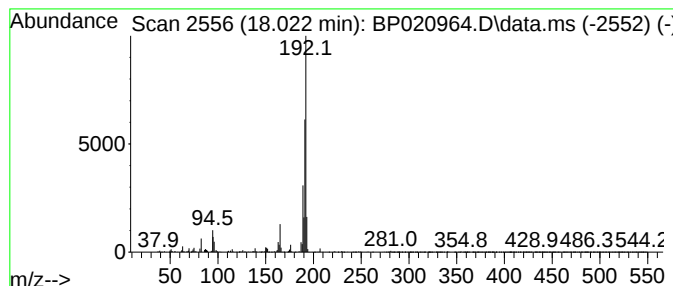
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	2.14 ng/ul	277071	Phenanthrene-d10	17.122

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



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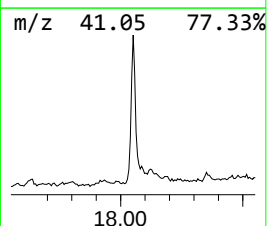
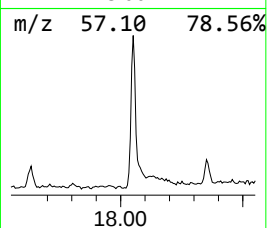
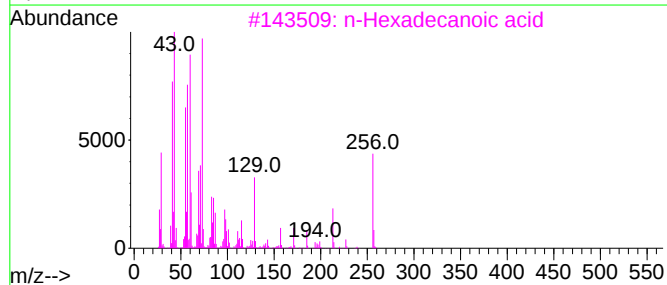
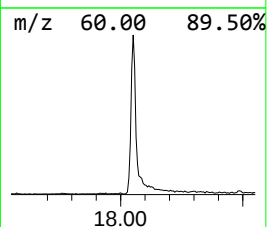
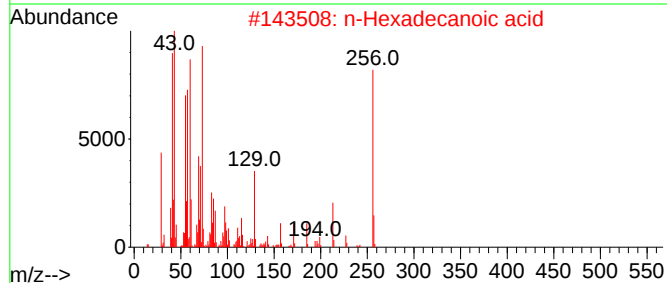
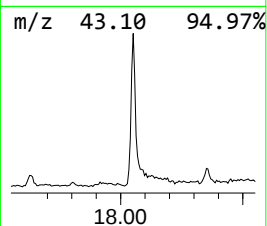
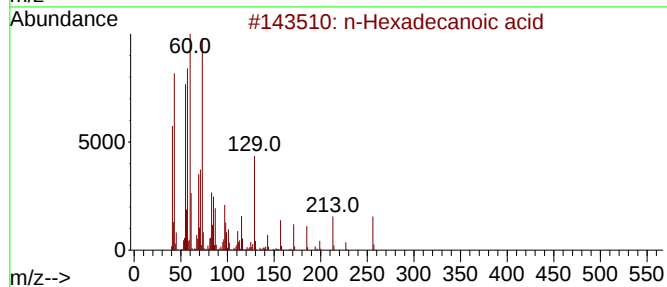
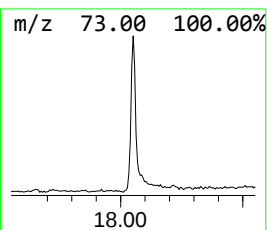
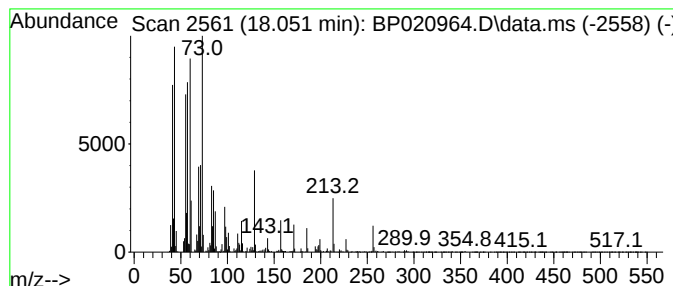
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	3.38 ng/ul	436646	Phenanthrene-d10	17.122

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
Operator : MA/JU
Sample : P3087-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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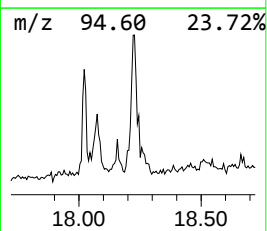
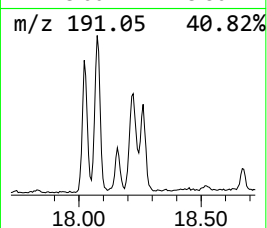
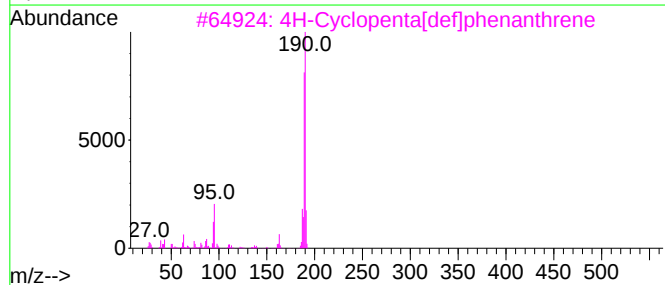
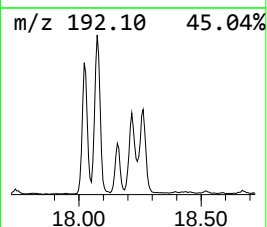
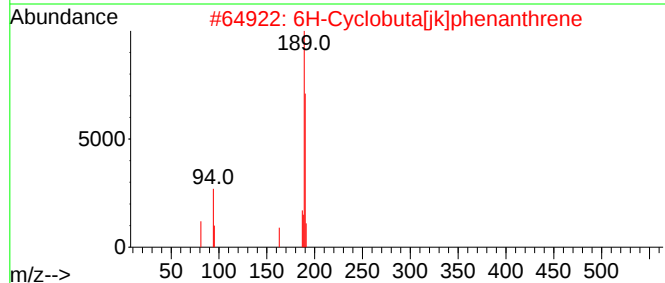
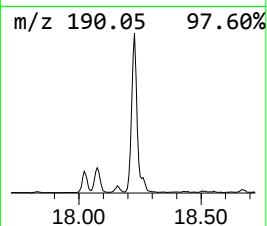
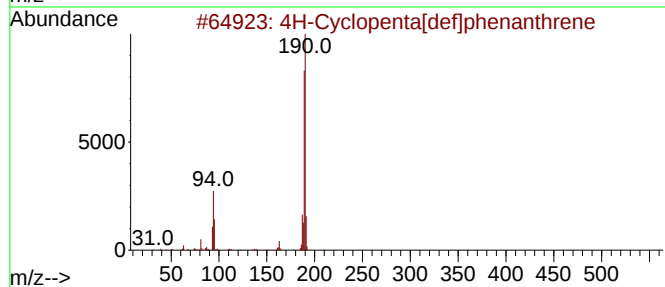
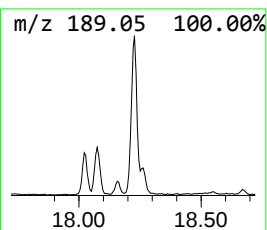
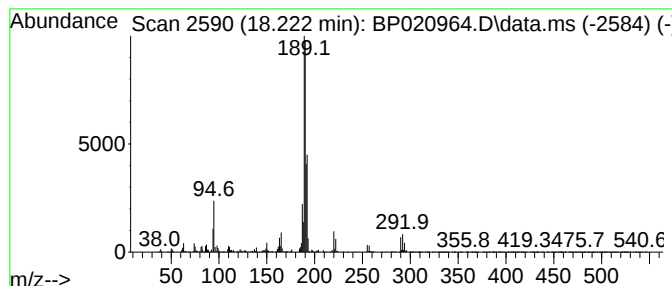
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	5.30 ng/ul	685144	Phenanthrene-d10	17.122

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5			4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
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Sample : P3087-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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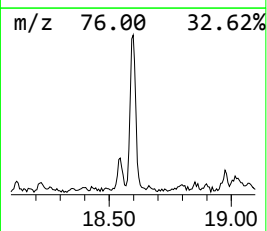
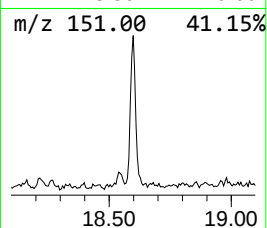
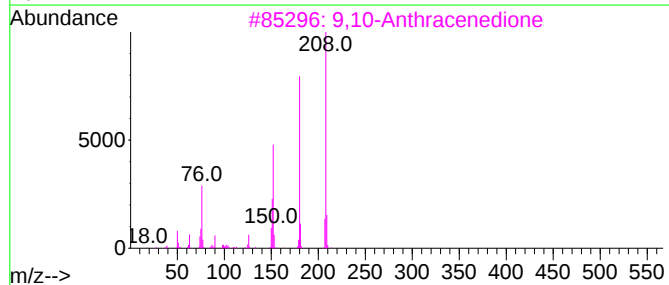
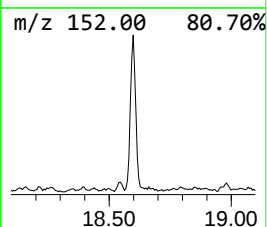
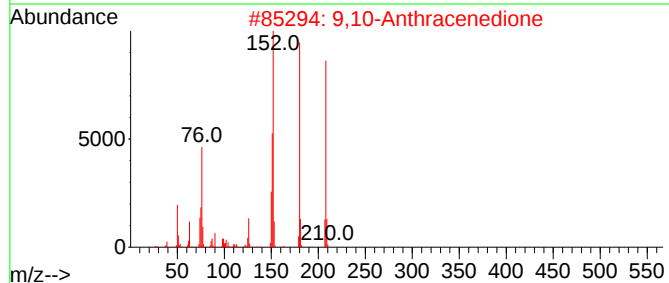
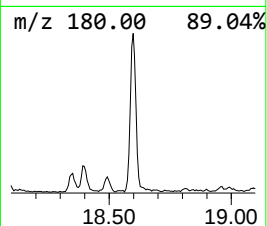
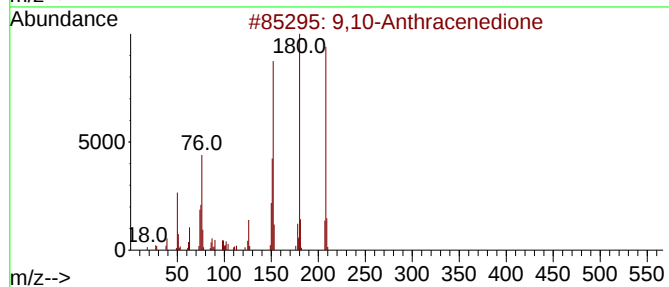
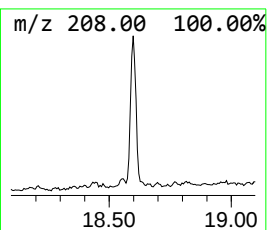
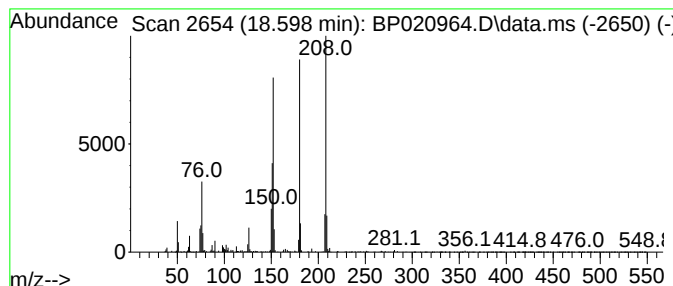
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.598	2.38 ng/ul	307719	Phenanthrene-d10	17.122

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
4		9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
Operator : MA/JU
Sample : P3087-01
Misc :
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Instrument :
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ClientSampleId :
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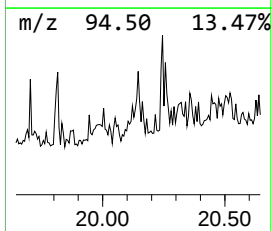
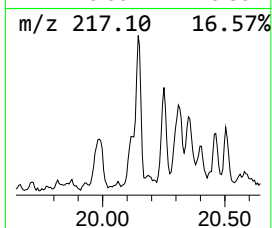
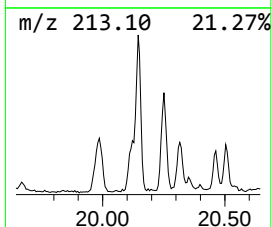
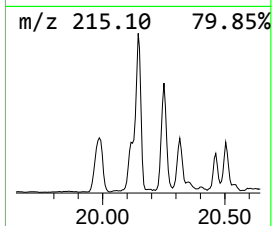
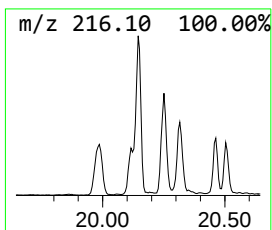
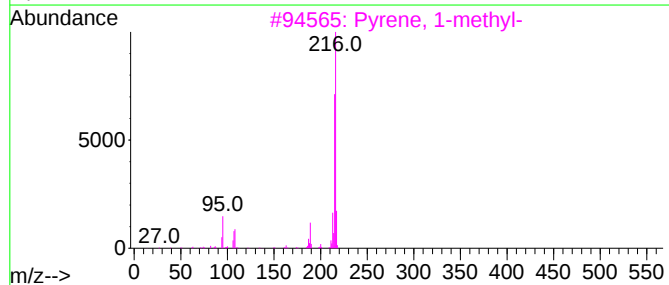
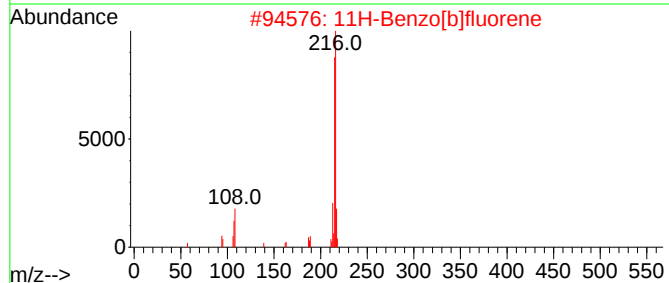
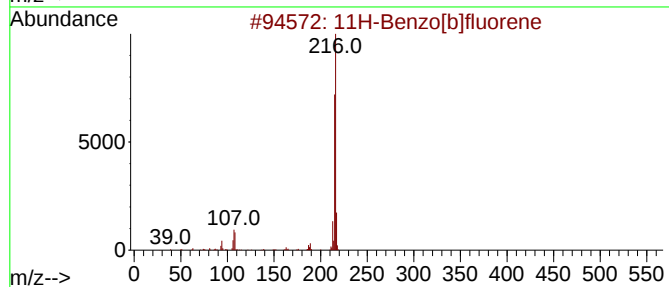
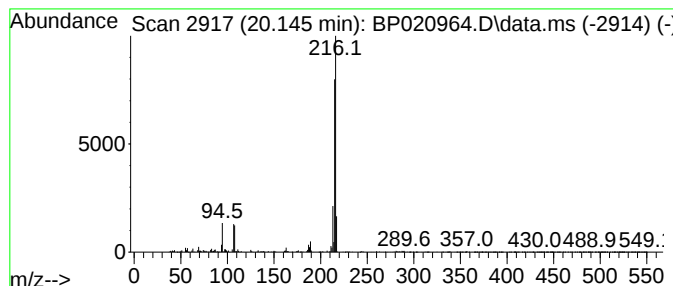
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.145	2.41 ng/ul	669603	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
Operator : MA/JU
Sample : P3087-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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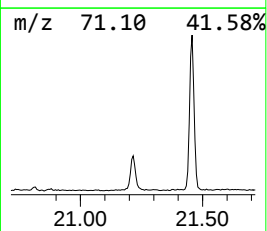
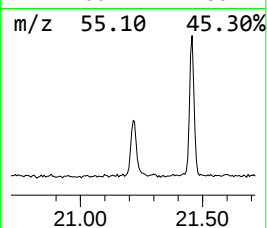
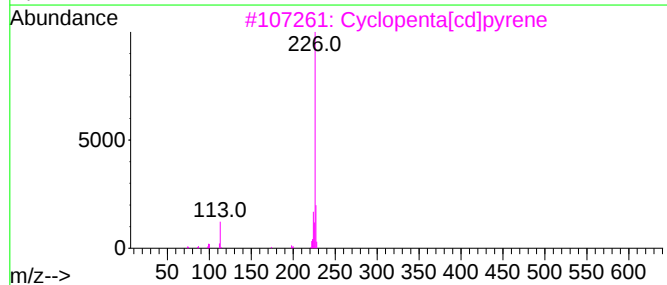
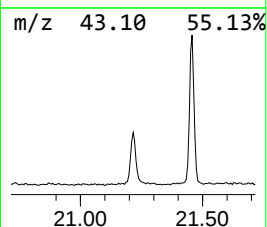
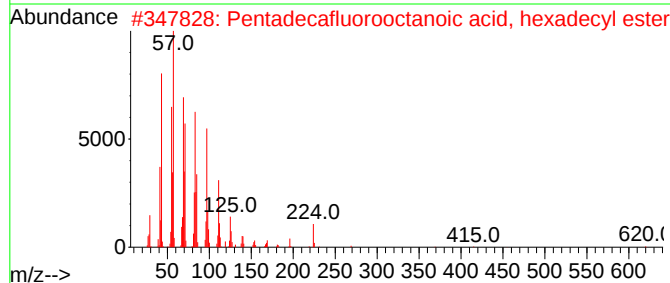
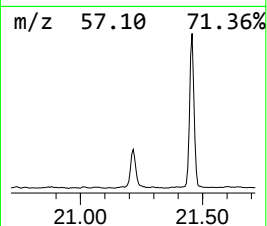
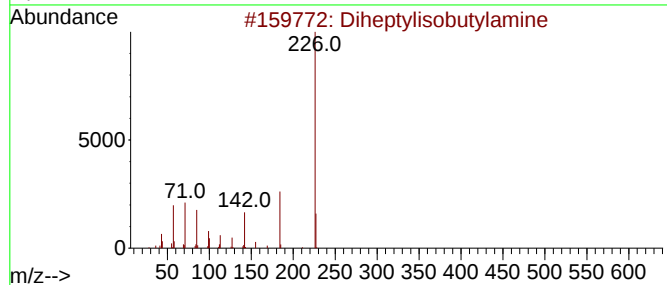
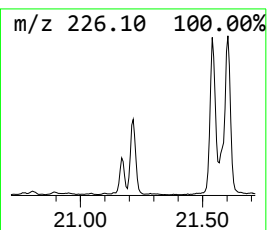
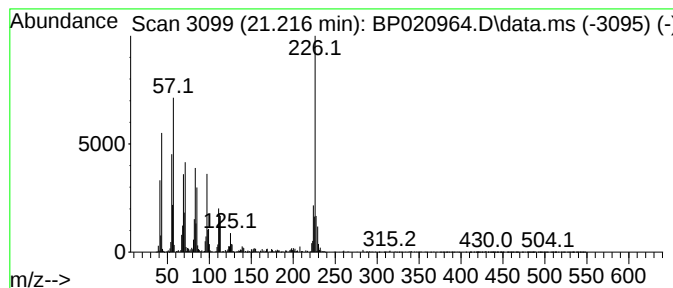
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.215	3.92 ng/ul	1087600	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diheptylisobutylamine	269	C18H39N	1000310-32-0	37
2			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	30
3			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	30
4			Butyramide, 2-phenyl-N-propyl-N...	345	C23H39NO	1000438-72-0	25
5			4-Amino-5-(4-chlorophenyl)-4H-1...	226	C8H7ClN4S	1010476-96-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
Operator : MA/JU
Sample : P3087-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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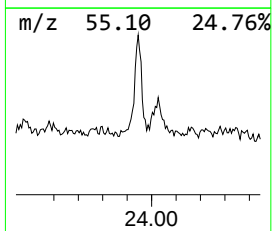
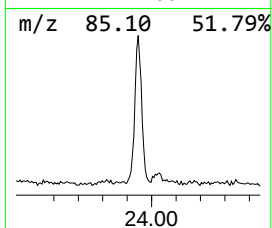
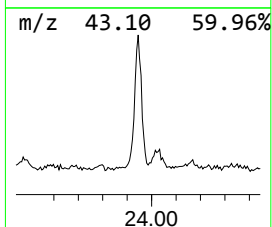
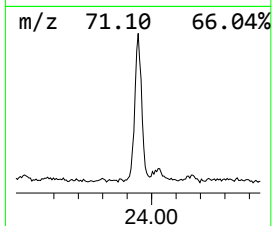
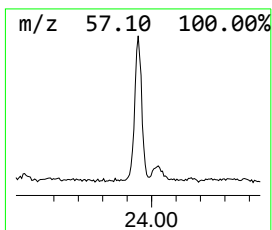
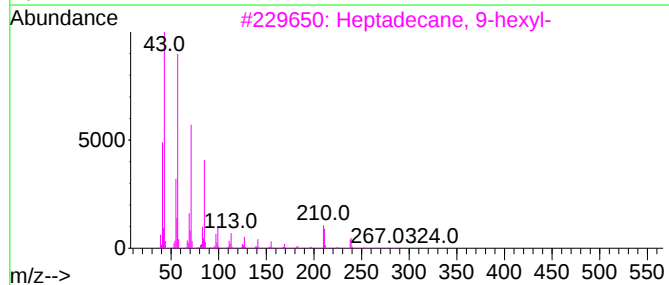
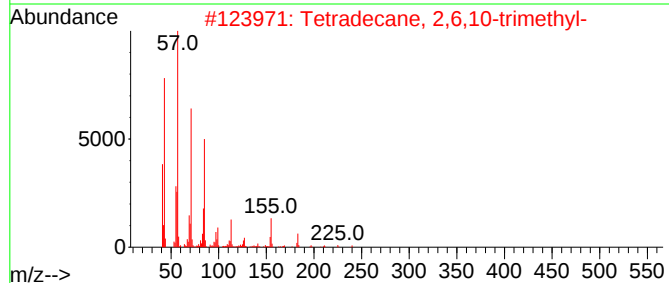
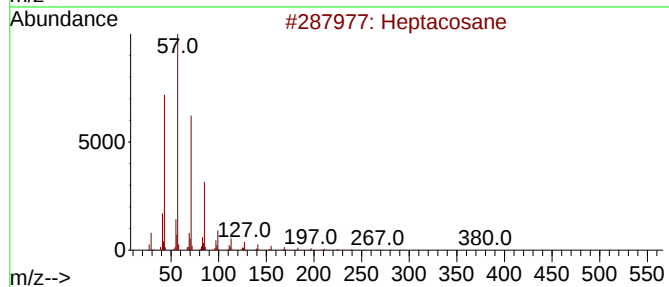
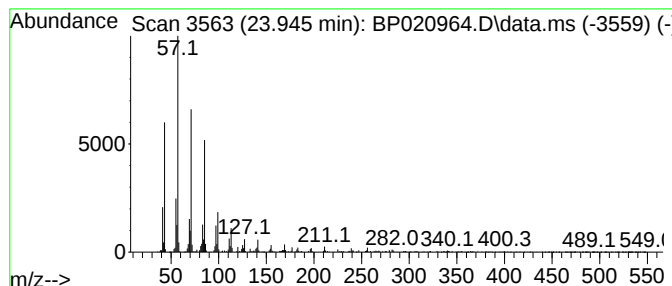
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.945	4.96 ng/ul	652980	Perylene-d12	24.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	94
2			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
3			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4			Hexacosane	366	C26H54	000630-01-3	91
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020964.D
Acq On : 15 Jul 2024 18:53
Operator : MA/JU
Sample : P3087-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Phenanthrene, 2...	18.022	2.1	ng/u1	277071	4	17.122	2585240	20.0
n-Hexadecanoic ...	18.051	3.4	ng/u1	436646	4	17.122	2585240	20.0
4H-Cyclopenta[d...	18.222	5.3	ng/u1	685144	4	17.122	2585240	20.0
9,10-Anthracene...	18.598	2.4	ng/u1	307719	4	17.122	2585240	20.0
11H-Benzo[b]flu...	20.145	2.4	ng/u1	669603	5	21.557	5546200	20.0
unknown-01	21.215	3.9	ng/u1	1087600	5	21.557	5546200	20.0
(DEL) Alkane: S...	23.945	5.0	ng/u1	652980	6	24.857	2634470	20.0