

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017350.D
 Acq On : 26 Sep 2023 08:18
 Operator : MA/JU
 Sample : 04472-10
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYK8

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-BP092023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017350.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.746	108	112	124	rBV	51421	93684	2.66%	0.189%
2	3.846	125	129	134	rVV	37757	50514	1.43%	0.102%
3	3.922	137	142	150	rVB	82636	121434	3.45%	0.245%
4	4.187	182	187	194	rVB4	27217	43507	1.24%	0.088%
5	4.428	223	228	234	rVB3	28487	42761	1.21%	0.086%
6	4.975	316	321	327	rBV	55238	84706	2.40%	0.171%
7	5.534	408	416	427	rBV	115560	218733	6.21%	0.442%
8	7.081	673	679	692	rBV	391090	689288	19.57%	1.391%
9	7.269	702	711	719	rBV	354699	545446	15.48%	1.101%
10	7.446	733	741	751	rVV	416055	680099	19.31%	1.373%
11	7.540	752	757	762	rVV	29115	44905	1.27%	0.091%
12	7.922	815	822	832	rBV	846031	1359037	38.58%	2.743%
13	8.640	937	944	952	rBV	339899	557530	15.83%	1.125%
14	8.775	962	967	975	rVB	142390	220791	6.27%	0.446%
15	9.122	1021	1026	1037	rVV	401727	659592	18.72%	1.331%
16	9.840	1140	1148	1155	rVV	324935	534753	15.18%	1.079%
17	9.981	1166	1172	1177	rBV	25003	48410	1.37%	0.098%
18	10.363	1230	1237	1248	rVV	491201	827125	23.48%	1.670%
19	10.640	1277	1284	1289	rVB	27829	45268	1.29%	0.091%
20	10.752	1295	1303	1309	rBV	1206747	2003821	56.88%	4.045%
21	10.804	1309	1312	1320	rVB2	55366	95415	2.71%	0.193%
22	10.916	1323	1331	1341	rBV	234047	435348	12.36%	0.879%
23	11.681	1456	1461	1466	rVV	29334	41174	1.17%	0.083%
24	12.093	1526	1531	1536	rBV2	39857	57801	1.64%	0.117%
25	12.404	1578	1584	1590	rVV	126208	200741	5.70%	0.405%
26	12.628	1616	1622	1627	rBV	73350	112504	3.19%	0.227%
27	13.340	1737	1743	1747	rBV	52840	75474	2.14%	0.152%
28	13.422	1752	1757	1766	rVV2	67891	129969	3.69%	0.262%
29	13.740	1804	1811	1812	rBV	56595	77614	2.20%	0.157%
30	13.857	1821	1831	1833	rBV4	44599	74437	2.11%	0.150%
31	13.898	1833	1838	1843	rVV	79383	147234	4.18%	0.297%
32	13.951	1843	1847	1850	rVV	79341	110152	3.13%	0.222%
33	13.998	1850	1855	1865	rVB	985251	1412868	40.11%	2.852%
34	14.140	1872	1879	1882	rBV	48905	76940	2.18%	0.155%
35	14.287	1895	1904	1911	rBV	1093033	1614828	45.84%	3.260%
36	14.416	1922	1926	1935	rVB	69341	88966	2.53%	0.180%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Title : SVOA CALIBRATION

37	14.587	1945	1955	1964	rVV	1877429	2732542	77.57%	5.516%
38	14.663	1964	1968	1973	rVB3	25579	43519	1.24%	0.088%
39	14.822	1990	1995	2008	rVV	178364	388872	11.04%	0.785%
40	14.969	2015	2020	2027	rVV3	69942	169196	4.80%	0.342%
41	15.045	2027	2033	2041	rVB2	65044	127829	3.63%	0.258%
42	15.187	2052	2057	2063	rBV2	45394	81547	2.31%	0.165%
43	15.245	2063	2067	2075	rVB	55353	80425	2.28%	0.162%
44	15.369	2077	2088	2098	rBV4	99128	236082	6.70%	0.477%
45	15.587	2114	2125	2135	rVV	1462637	2224129	63.14%	4.490%
46	15.663	2135	2138	2143	rVB2	31703	43743	1.24%	0.088%
47	15.716	2143	2147	2152	rBV	138926	195186	5.54%	0.394%
48	15.769	2152	2156	2162	rVB5	38365	68532	1.95%	0.138%
49	15.928	2179	2183	2188	rVB2	69780	104453	2.97%	0.211%
50	16.022	2193	2199	2204	rBV	213042	276242	7.84%	0.558%
51	16.075	2204	2208	2216	rBV	45684	95265	2.70%	0.192%
52	16.169	2221	2224	2230	rVB2	41731	55140	1.57%	0.111%
53	16.245	2232	2237	2243	rVV	127524	229679	6.52%	0.464%
54	16.292	2243	2245	2251	rVB4	32026	48807	1.39%	0.099%
55	16.645	2297	2305	2309	rVB3	55379	122335	3.47%	0.247%
56	16.692	2309	2313	2317	rBV2	72152	111705	3.17%	0.225%
57	16.869	2336	2343	2346	rBV2	76575	143950	4.09%	0.291%
58	16.904	2346	2349	2353	rVB2	58202	76592	2.17%	0.155%
59	17.010	2360	2367	2370	rBV4	66385	156974	4.46%	0.317%
60	17.039	2370	2372	2375	rVV	69406	83305	2.36%	0.168%
61	17.075	2375	2378	2384	rVB5	64491	113296	3.22%	0.229%
62	17.175	2390	2395	2409	rVB4	54801	152306	4.32%	0.307%
63	17.357	2420	2426	2431	rBV	2199784	3200435	90.85%	6.460%
64	17.404	2431	2434	2438	rVB	403507	507365	14.40%	1.024%
65	17.457	2438	2443	2448	rBV	1471635	2061425	58.52%	4.161%
66	17.628	2467	2472	2476	rVB3	57403	91576	2.60%	0.185%
67	17.675	2476	2480	2486	rVB2	98675	154857	4.40%	0.313%
68	17.786	2488	2499	2508	rBV7	167471	457590	12.99%	0.924%
69	17.886	2511	2516	2517	rBV4	38682	59750	1.70%	0.121%
70	17.939	2521	2525	2529	rVV2	92752	127945	3.63%	0.258%
71	18.151	2557	2561	2566	rBV	321890	442626	12.56%	0.893%
72	18.210	2566	2571	2576	rVV2	496548	1077098	30.58%	2.174%
73	18.269	2577	2581	2586	rVB	427453	618037	17.54%	1.248%
74	18.345	2591	2594	2597	rVV2	95256	122020	3.46%	0.246%
75	18.404	2600	2604	2608	rVV	262034	383224	10.88%	0.774%
76	18.445	2608	2611	2621	rVB2	216871	400362	11.36%	0.808%

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77	18.528	2621	2625	2627	rBV4	70069	90714	2.58%	0.183%
78	18.704	2651	2655	2661	rBV2	157131	237906	6.75%	0.480%
79	18.939	2692	2695	2699	rBV2	93781	111283	3.16%	0.225%
80	18.986	2700	2703	2706	rVB	159048	174922	4.97%	0.353%
81	19.028	2706	2710	2715	rBV3	144943	185455	5.26%	0.374%
82	19.104	2715	2723	2727	rBV2	386022	706080	20.04%	1.425%
83	19.151	2728	2731	2735	rBV	120718	173639	4.93%	0.351%
84	19.192	2735	2738	2742	rVB	158341	173201	4.92%	0.350%
85	19.233	2742	2745	2746	rBV2	133539	136672	3.88%	0.276%
86	19.333	2759	2762	2763	rBV2	63936	74089	2.10%	0.150%
87	19.369	2764	2768	2772	rVB2	261939	410196	11.64%	0.828%
88	19.433	2775	2779	2792	rVB3	160389	477854	13.56%	0.965%
89	19.639	2809	2814	2818	rVB2	176559	256943	7.29%	0.519%
90	19.704	2819	2825	2835	rBV	2252598	3395886	96.40%	6.855%
91	19.786	2835	2839	2844	rVB2	184025	287944	8.17%	0.581%
92	20.110	2891	2894	2897	rVB3	115828	122276	3.47%	0.247%
93	20.163	2900	2903	2908	rVV4	203778	281975	8.00%	0.569%
94	20.651	2983	2986	2989	rVB	235088	253350	7.19%	0.511%
95	21.110	3060	3064	3066	rBV	342809	406367	11.54%	0.820%
96	21.463	3119	3124	3129	rBV	2807936	3522778	100.00%	7.111%
97	21.551	3136	3139	3144	rVB2	279740	300365	8.53%	0.606%
98	22.022	3216	3219	3223	rBV	275637	317552	9.01%	0.641%
99	23.810	3518	3523	3530	rVB	1305272	2511876	71.30%	5.070%
100	23.980	3545	3552	3560	rBV	1664500	3441634	97.70%	6.947%

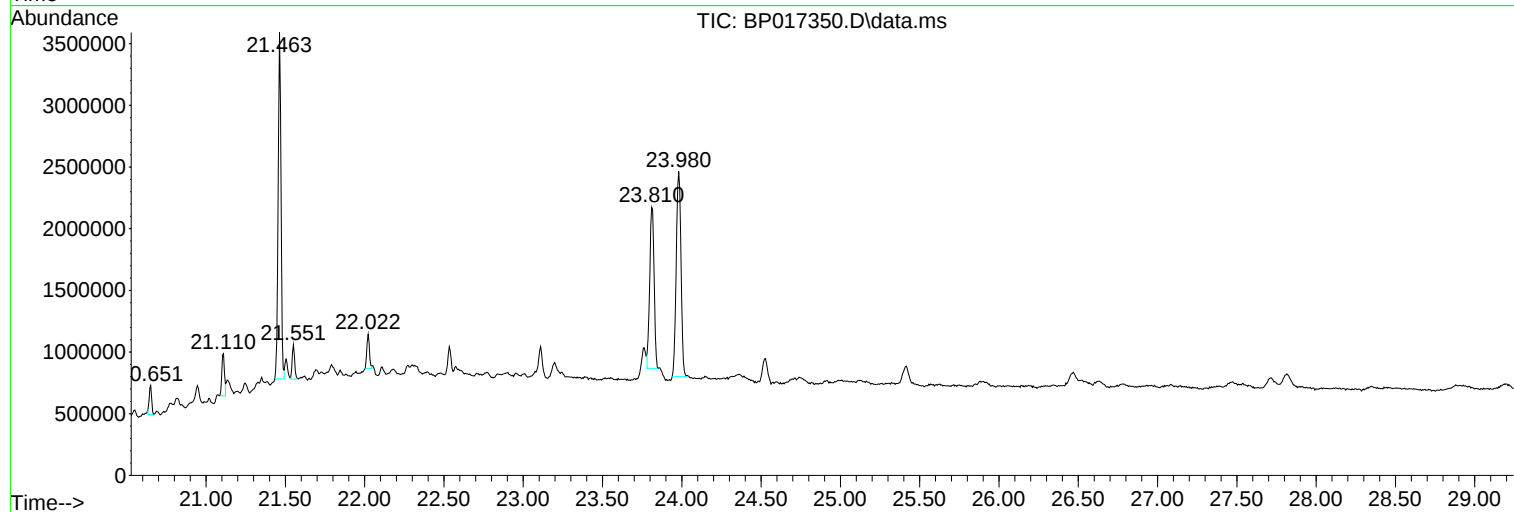
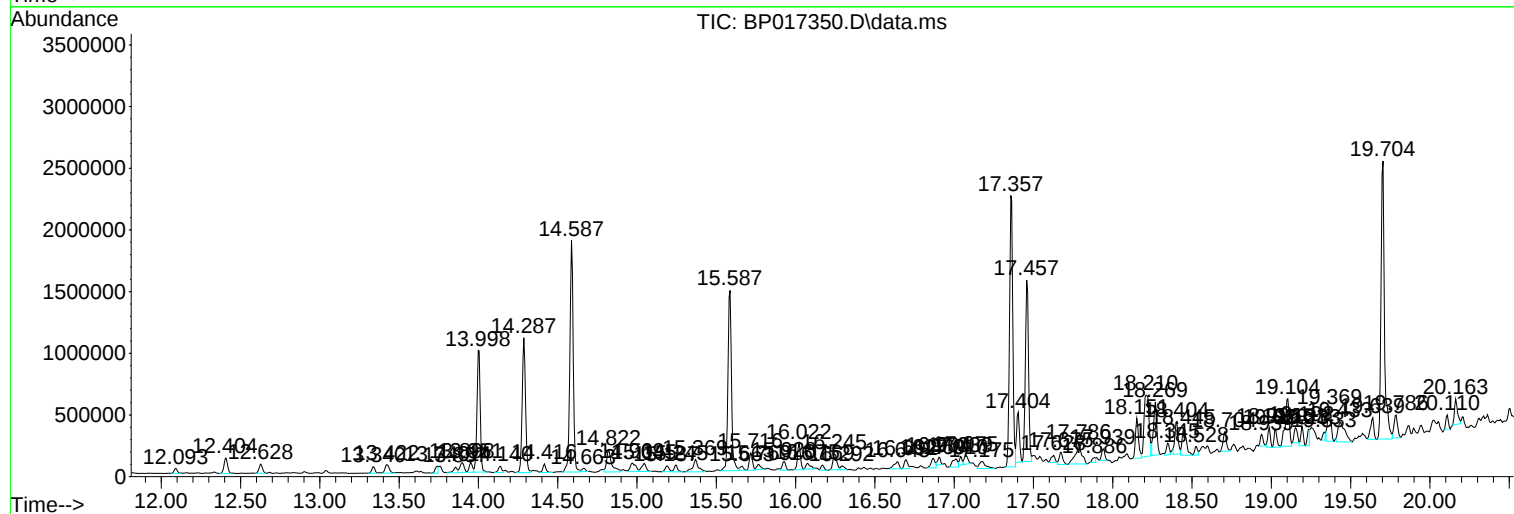
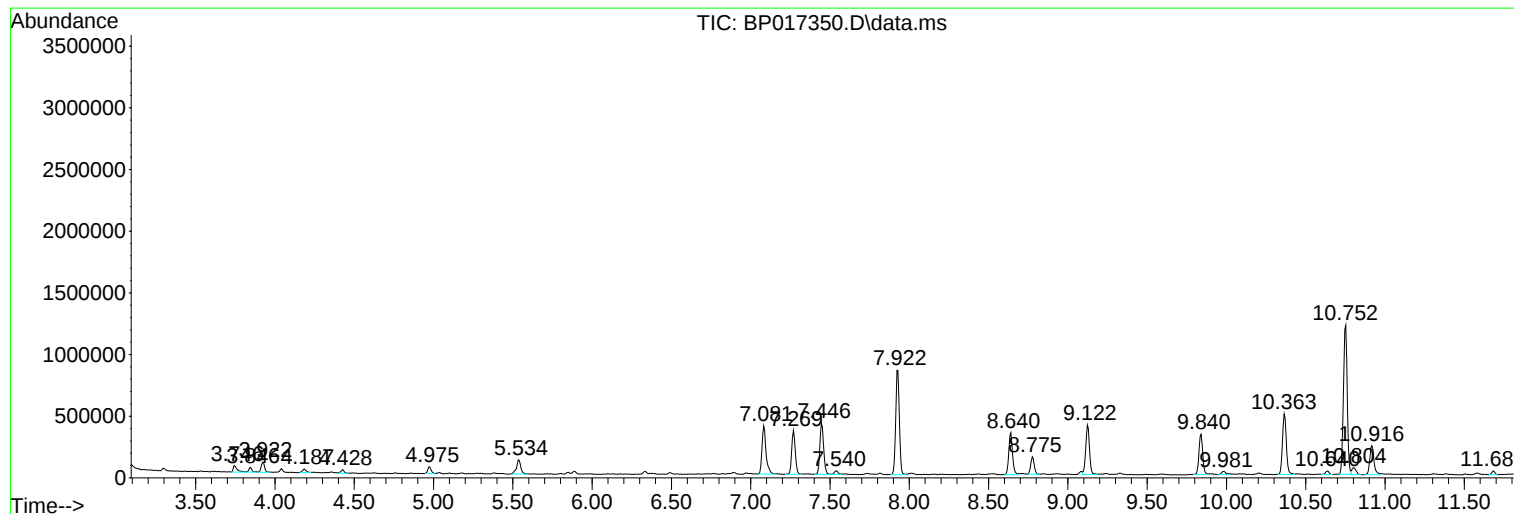
Sum of corrected areas: 49539787

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



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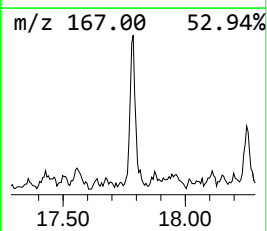
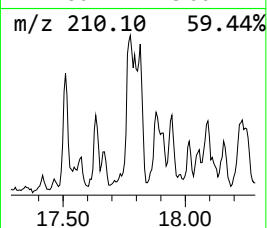
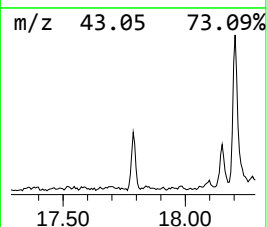
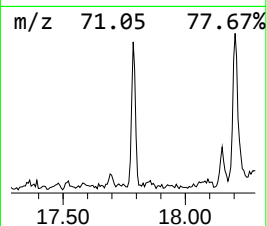
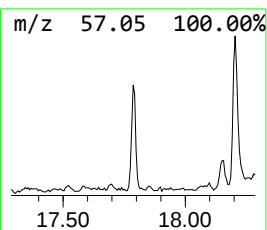
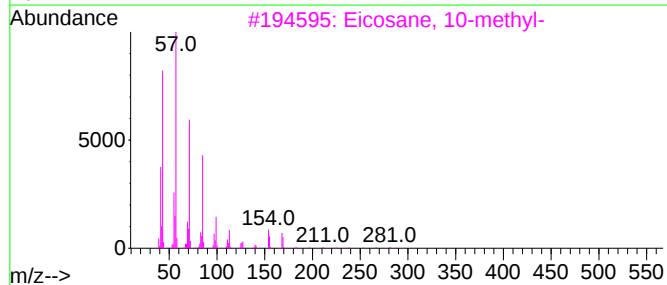
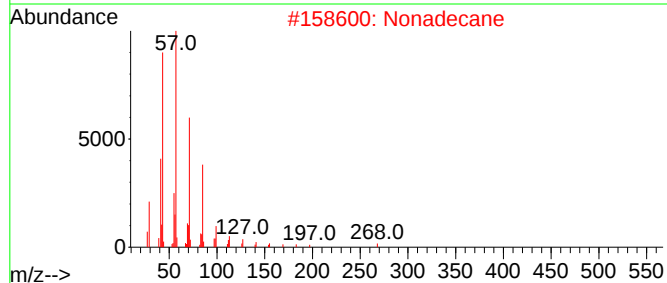
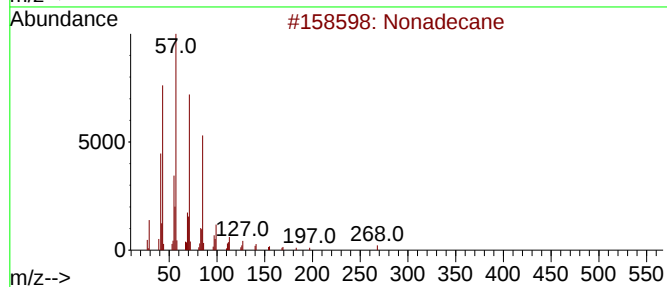
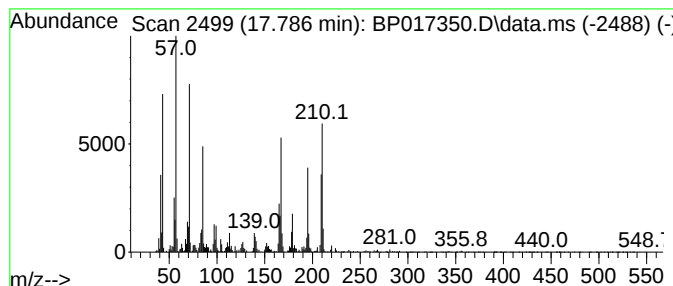
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.787	2.86 ng/ul	457590	Phenanthrene-d10	17.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	86
2		Nonadecane	268	C19H40	000629-92-5	50
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	45
4		Heneicosane	296	C21H44	000629-94-7	42
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	42



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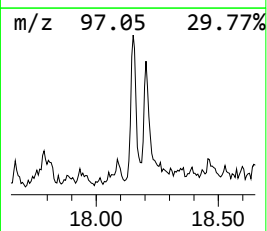
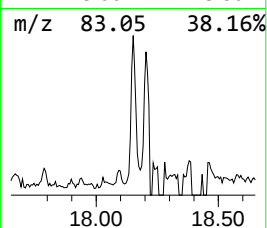
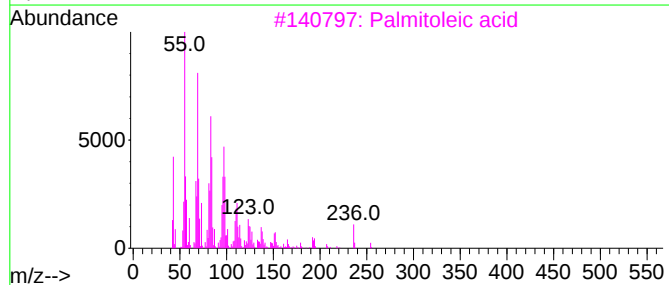
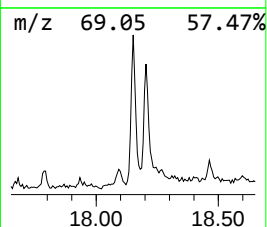
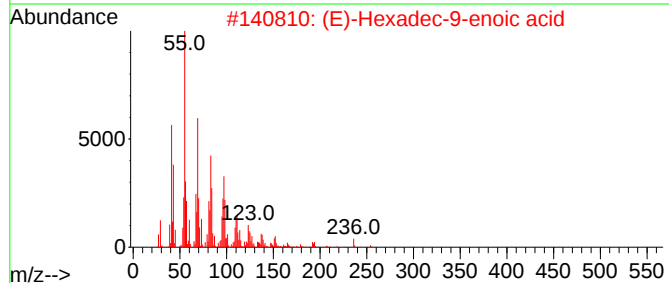
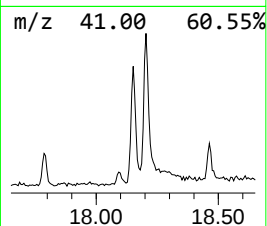
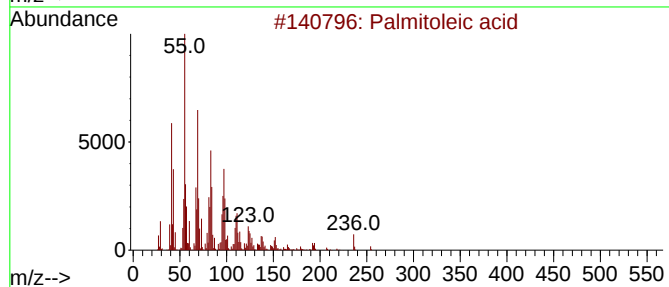
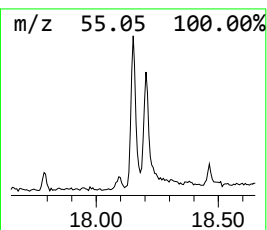
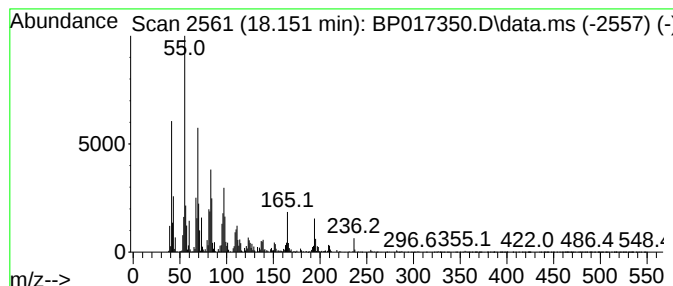
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Palmitoleic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	2.77 ng/ul	442626	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	99
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
3			Palmitoleic acid	254	C16H30O2	000373-49-9	97
4			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	96
5			13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29BO	1000156-41-7	93



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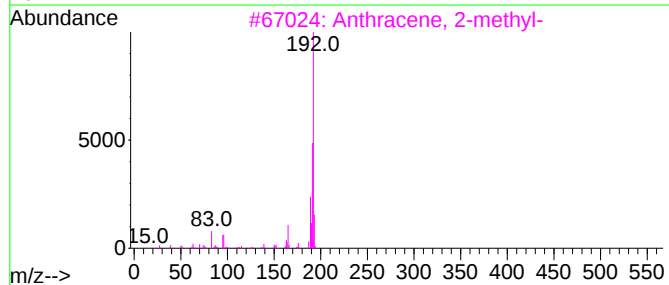
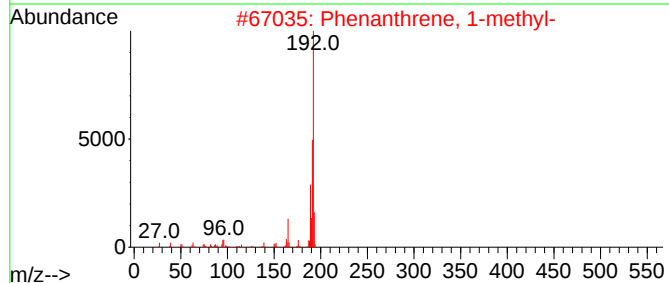
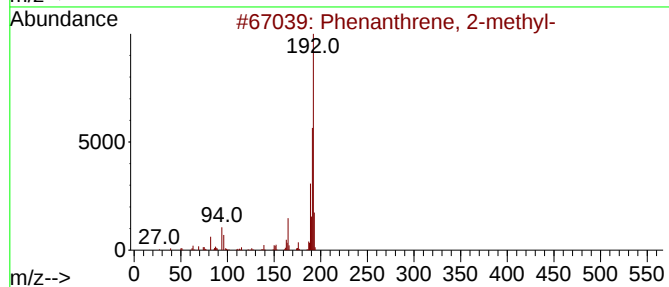
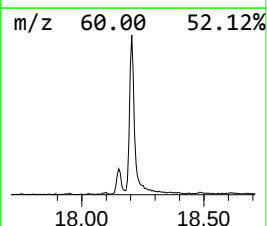
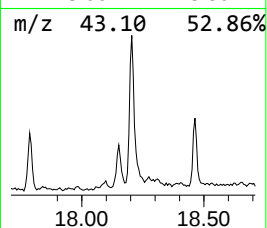
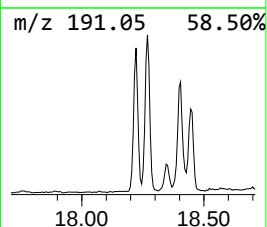
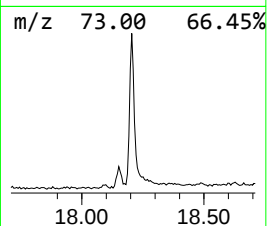
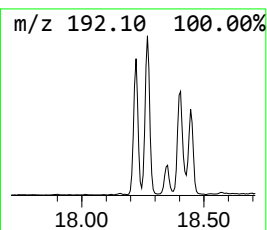
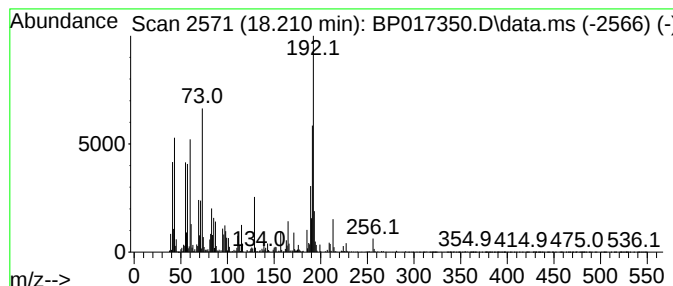
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	6.73 ng/ul	1077100	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYK8

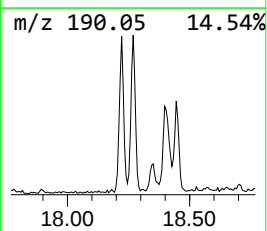
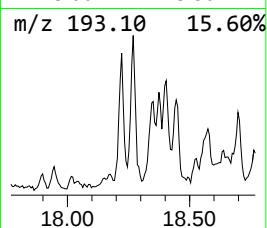
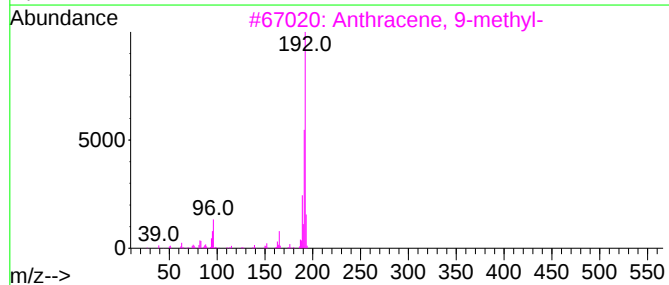
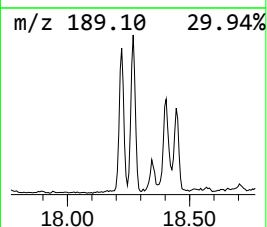
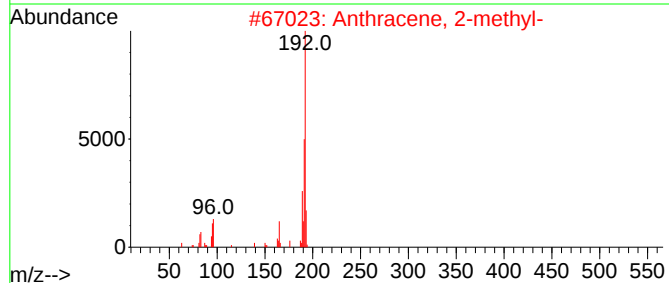
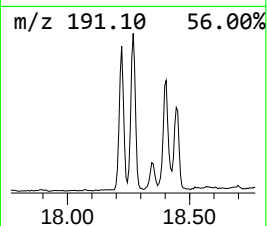
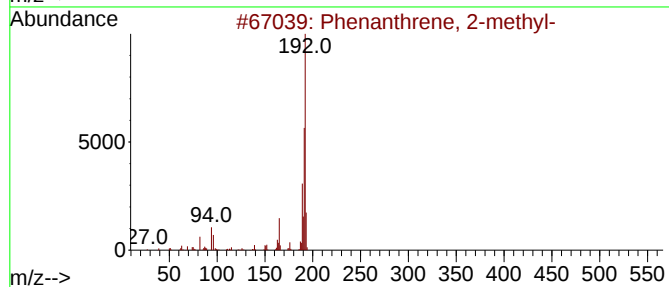
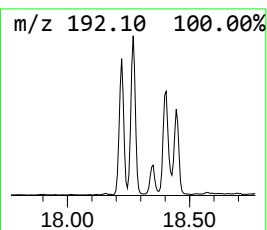
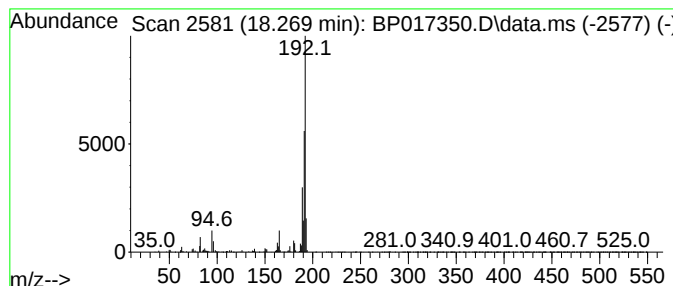
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.269	3.86 ng/ul	618037	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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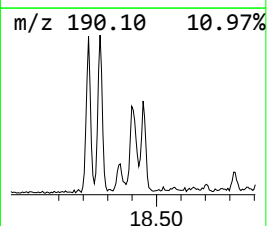
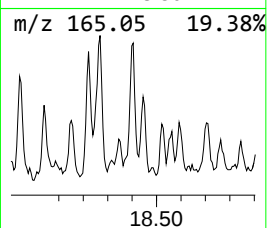
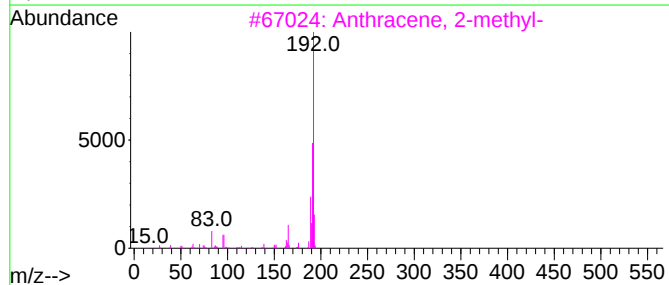
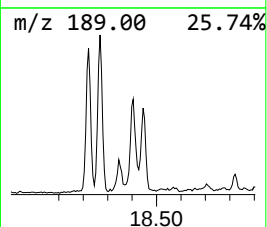
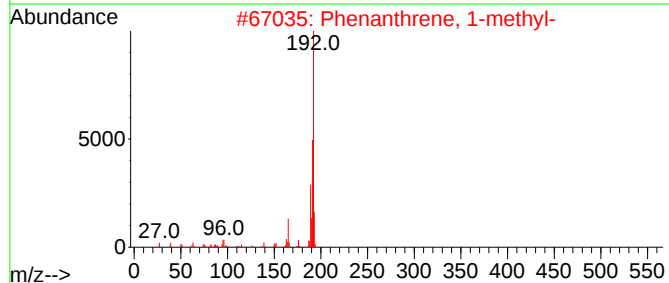
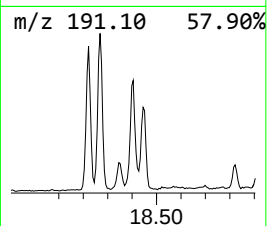
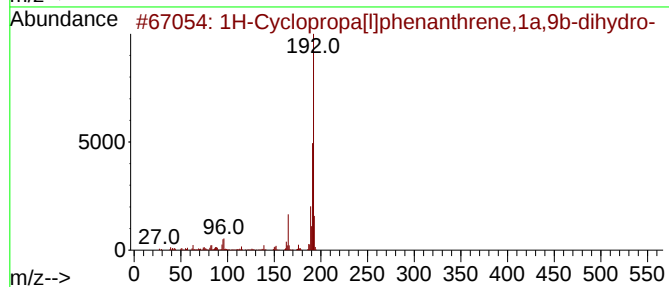
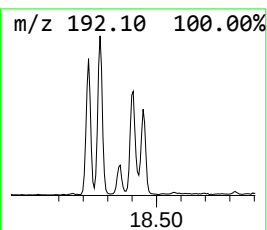
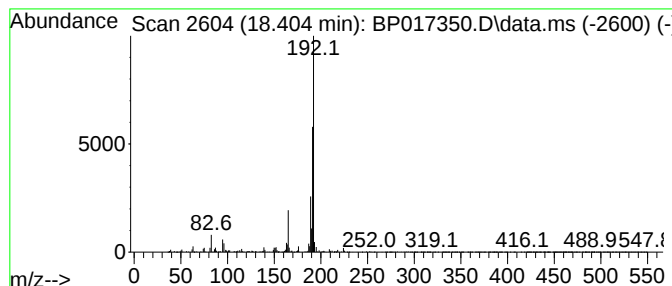
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	2.39 ng/ul	383224	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
4			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	87
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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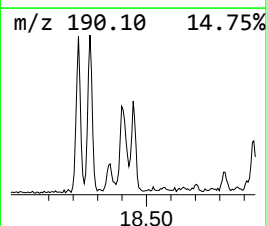
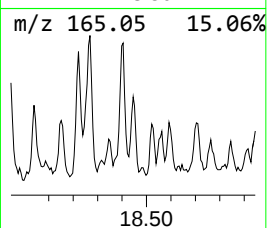
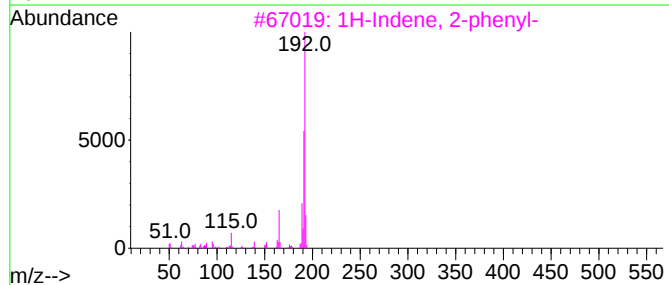
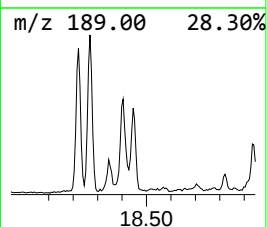
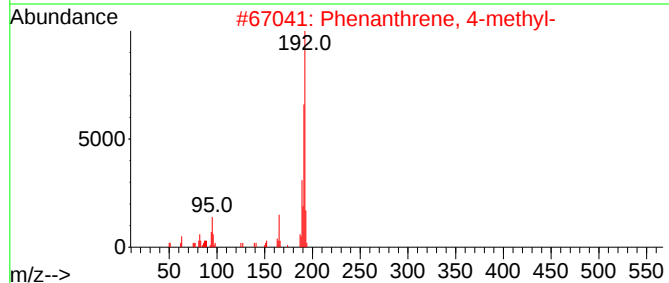
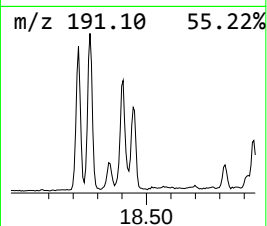
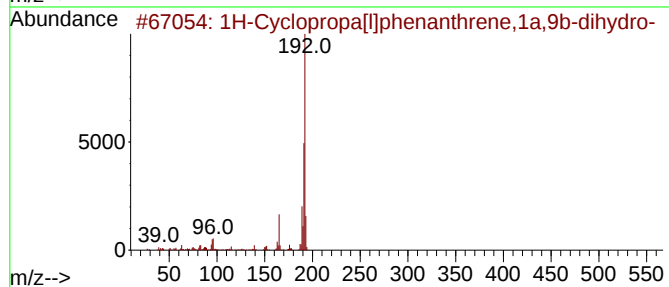
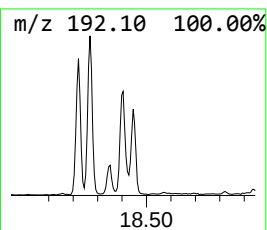
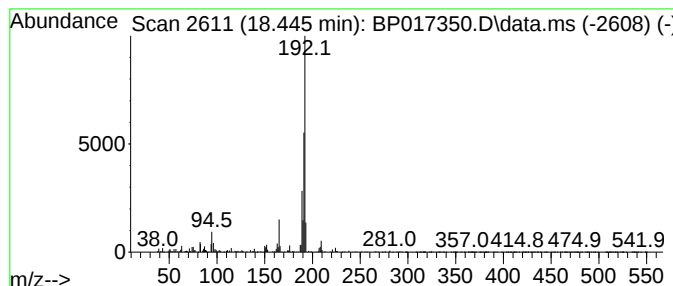
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenanthrene, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.445	2.50 ng/ul	400362	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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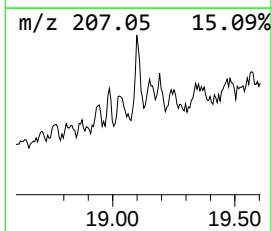
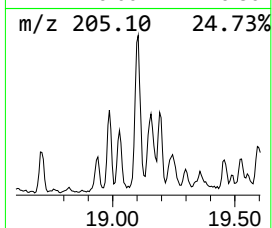
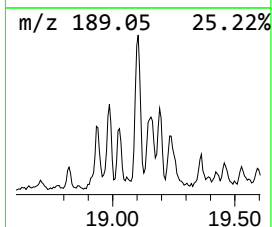
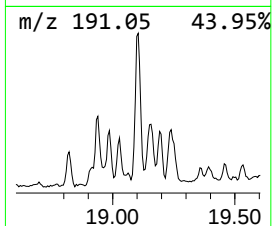
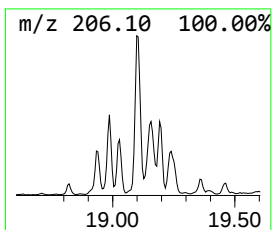
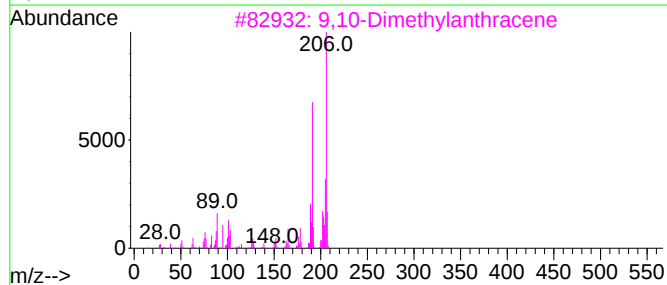
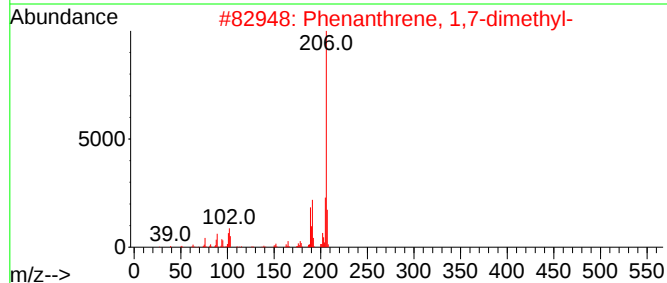
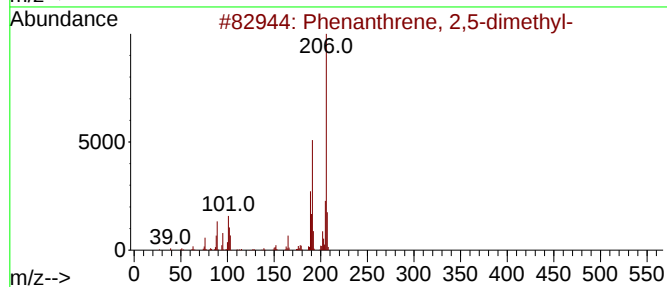
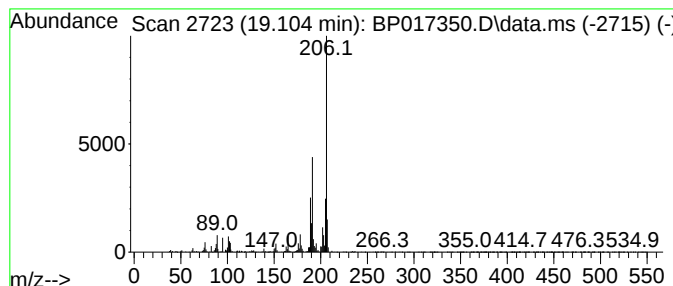
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.104	4.41 ng/ul	706080	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3			9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4			9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
5			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

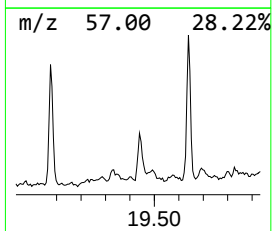
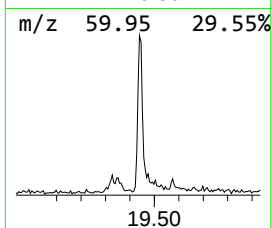
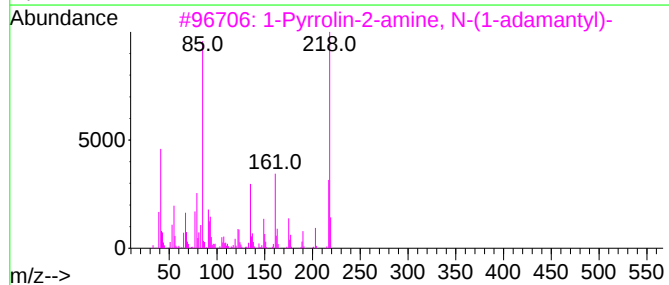
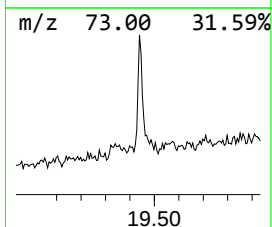
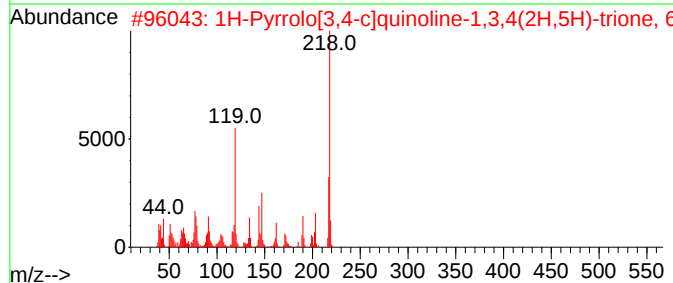
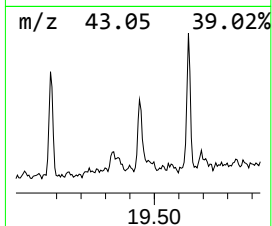
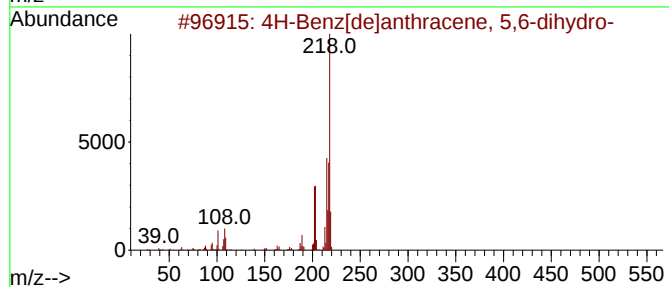
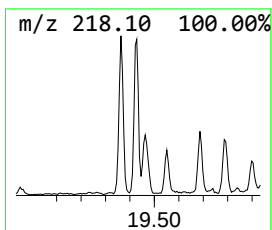
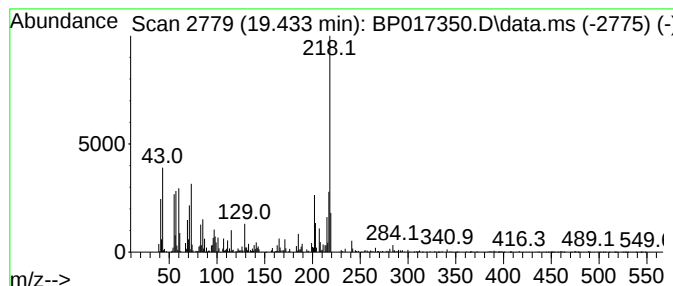
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 4H-Benz[de]anthracene, 5,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.433	2.71 ng/ul	477854	Chrysene-d12		21.463	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Benz[de]anthracene, 5,6-dihydro-		218	C17H14	004389-09-7	58
2	1H-Pyrrolo[3,4-c]quinoline-1,3,4...		218	C11H10N2O3	181021-85-2	47
3	1-Pyrrolin-2-amine, N-(1-adamant...		218	C14H22N2	300695-00-5	43
4	3,4-Dihydro-2-bromo-4,4-dimethyl...		312	C18H17Br	071161-44-9	43
5	Naphthalene, 1-(4-methylphenyl)-		218	C17H14	027331-34-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

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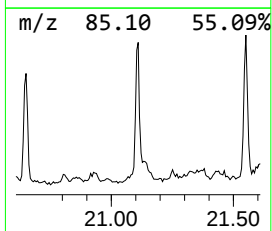
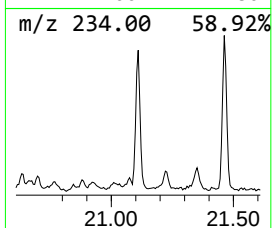
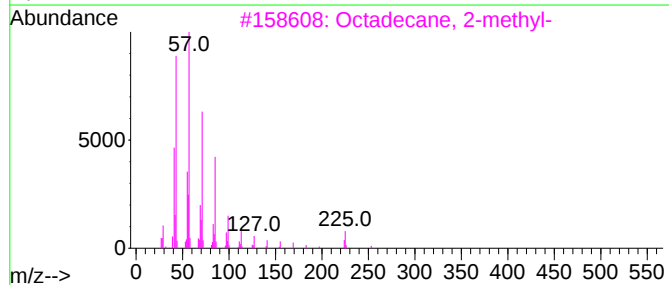
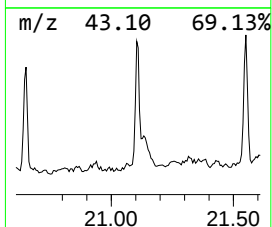
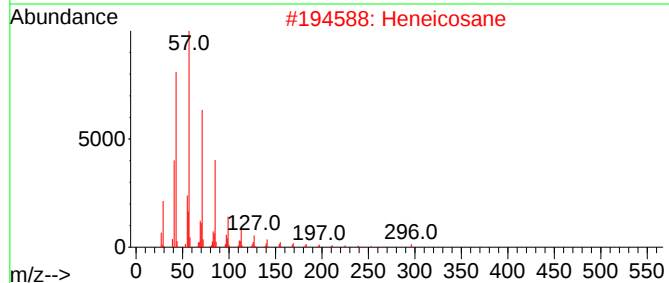
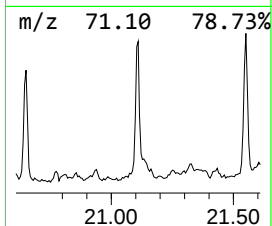
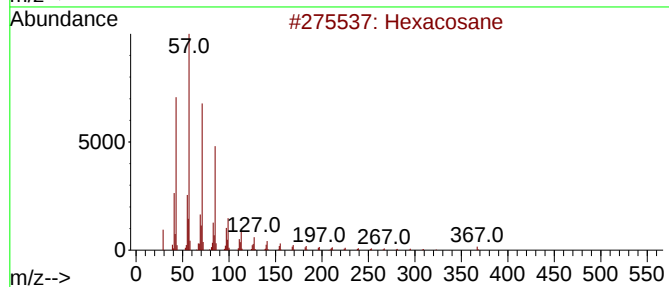
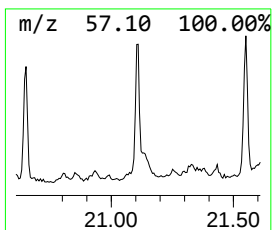
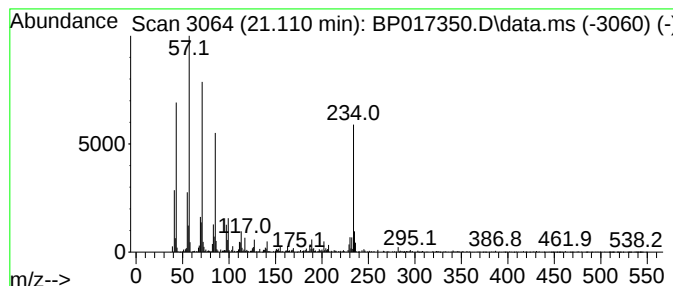
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	2.31 ng/ul	406367	Chrysene-d12	21.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	70
2		Heneicosane	296	C21H44	000629-94-7	58
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	58
4		Heptacosane	380	C27H56	000593-49-7	58
5		Nonadecane	268	C19H40	000629-92-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017350.D
Acq On : 26 Sep 2023 08:18
Operator : MA/JU
Sample : 04472-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.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
(DEL) Alkane: S...	17.787	2.9	ng/ul	457590	4	17.357	3200440	20.0
Palmitoleic acid	18.151	2.8	ng/ul	442626	4	17.357	3200440	20.0
Phenanthrene, 2...	18.210	6.7	ng/ul	1077100	4	17.357	3200440	20.0
Anthracene, 2-m...	18.269	3.9	ng/ul	618037	4	17.357	3200440	20.0
1H-Cyclopropa[1...	18.404	2.4	ng/ul	383224	4	17.357	3200440	20.0
Phenanthrene, 4...	18.445	2.5	ng/ul	400362	4	17.357	3200440	20.0
Phenanthrene, 2...	19.104	4.4	ng/ul	706080	4	17.357	3200440	20.0
4H-Benz[de]anth...	19.433	2.7	ng/ul	477854	5	21.463	3522780	20.0
(DEL) Alkane: S...	21.110	2.3	ng/ul	406367	5	21.463	3522780	20.0