

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043763.D
 Acq On : 05 Jan 2024 16:49
 Operator : MA/JU
 Sample : 05952-07ME
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF90ME

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_M\Methods\SFAM-EPA-BM121523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043763.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.357	26	29	40	rVB	53301	78637	0.11%	0.041%
2	7.128	662	670	684	rBV	804212	1357491	1.84%	0.701%
3	7.292	692	698	708	rBV	641377	1014223	1.38%	0.523%
4	7.481	722	730	739	rBV	803171	1311800	1.78%	0.677%
5	7.951	804	810	817	rBV	1111244	1726496	2.34%	0.891%
6	8.016	817	821	828	rVB	43500	76299	0.10%	0.039%
7	8.675	926	933	942	rBV	941709	1614754	2.19%	0.833%
8	8.792	947	953	963	rVB	58504	112140	0.15%	0.058%
9	9.128	1004	1010	1021	rBV	723122	1250523	1.70%	0.645%
10	9.851	1125	1133	1141	rBV	621800	1074747	1.46%	0.555%
11	10.392	1217	1225	1245	rBV	882087	1659993	2.25%	0.857%
12	10.763	1276	1288	1297	rBV	1371033	2378348	3.22%	1.227%
13	10.922	1304	1315	1331	rBV	361936	826640	1.12%	0.427%
14	12.175	1518	1528	1531	rBV8	34526	90371	0.12%	0.047%
15	12.822	1624	1638	1646	rBV2	93735	192270	0.26%	0.099%
16	13.645	1772	1778	1787	rVB10	28615	81761	0.11%	0.042%
17	13.769	1793	1799	1807	rBV3	50863	132258	0.18%	0.068%
18	13.921	1819	1825	1830	rBV	116029	222036	0.30%	0.115%
19	13.974	1830	1834	1836	rVV	64722	94883	0.13%	0.049%
20	14.016	1836	1841	1847	rVV	1640575	2300112	3.12%	1.187%
21	14.292	1882	1888	1901	rBV	2009210	3056505	4.14%	1.577%
22	14.486	1915	1921	1930	rBV	248643	462557	0.63%	0.239%
23	14.598	1934	1940	1946	rVV	1986862	2934880	3.98%	1.515%
24	14.657	1946	1950	1956	rBV5	87753	181392	0.25%	0.094%
25	14.716	1957	1960	1966	rVB5	88643	141725	0.19%	0.073%
26	14.833	1970	1980	1986	rBV2	754442	1506422	2.04%	0.777%
27	14.992	2001	2007	2014	rVB3	256520	460439	0.62%	0.238%
28	15.068	2016	2020	2024	rBV	227563	300986	0.41%	0.155%
29	15.145	2029	2033	2036	rVV4	104814	129913	0.18%	0.067%
30	15.227	2037	2047	2051	rVV2	904490	1886890	2.56%	0.974%
31	15.268	2051	2054	2058	rVB	196410	241790	0.33%	0.125%
32	15.386	2068	2074	2076	rBV2	128621	235531	0.32%	0.122%
33	15.433	2076	2082	2088	rVB3	414447	830436	1.13%	0.429%
34	15.592	2103	2109	2114	rBV	2491177	3572229	4.84%	1.843%
35	15.639	2114	2117	2121	rVB4	165392	232342	0.32%	0.120%
36	15.704	2122	2128	2129	rBV2	419369	671368	0.91%	0.346%

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Integration Parameters: LSCINT.P

Integrator: RTE

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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_M\Methods\SFAM-EPA-BM121523.MA.M
 Title : SVOA CALIBRATION

37	15.727	2129	2132	2142	rVB2	868724	1276357	1.73%	0.659%
38	15.845	2145	2152	2161	rVB2	579822	1291548	1.75%	0.666%
39	15.939	2163	2168	2174	rVB3	274687	574044	0.78%	0.296%
40	16.004	2175	2179	2182	rBV5	138343	244707	0.33%	0.126%
41	16.057	2185	2188	2191	rBV4	187884	225314	0.31%	0.116%
42	16.092	2191	2194	2200	rVV4	110896	205074	0.28%	0.106%
43	16.321	2225	2233	2238	rBV2	2056303	4474629	6.07%	2.309%
44	16.704	2294	2298	2303	rVB2	871879	1204829	1.63%	0.622%
45	16.880	2325	2328	2331	rBV4	333948	489227	0.66%	0.252%
46	17.015	2343	2351	2359	rBV	39802705	73748351	100.00%	38.057%
47	17.098	2361	2365	2368	rVV	2244370	2871723	3.89%	1.482%
48	17.145	2368	2373	2384	rVB2	2977062	6242351	8.46%	3.221%
49	17.351	2403	2408	2412	rBV	2507740	3853189	5.22%	1.988%
50	17.392	2412	2415	2419	rVV	1399187	2000528	2.71%	1.032%
51	17.451	2420	2425	2433	rVB	2793243	4345925	5.89%	2.243%
52	17.757	2474	2477	2482	rVB	953451	1410862	1.91%	0.728%
53	17.839	2488	2491	2498	rVB	2881270	3967668	5.38%	2.047%
54	17.933	2503	2507	2515	rVB3	1350477	2595650	3.52%	1.339%
55	18.233	2553	2558	2562	rBV2	2050838	4770310	6.47%	2.462%
56	18.274	2562	2565	2570	rVB2	3050104	4533000	6.15%	2.339%
57	18.415	2586	2589	2593	rVB	962657	1186418	1.61%	0.612%
58	18.539	2606	2610	2615	rBV2	2660009	4198867	5.69%	2.167%
59	18.621	2621	2624	2634	rVB4	1414166	2309940	3.13%	1.192%
60	18.974	2681	2684	2690	rBV5	797460	1525536	2.07%	0.787%
61	19.021	2690	2692	2696	rVB	987659	1041792	1.41%	0.538%
62	19.062	2696	2699	2708	rVB3	1438057	3104911	4.21%	1.602%
63	19.145	2709	2713	2715	rBV2	1867251	2643919	3.59%	1.364%
64	19.239	2726	2729	2733	rVB3	833870	1010203	1.37%	0.521%
65	19.462	2764	2767	2771	rBV3	853331	1406119	1.91%	0.726%
66	19.745	2811	2815	2819	rBV2	3796694	4943291	6.70%	2.551%
67	20.280	2904	2906	2910	rVB	643861	680193	0.92%	0.351%
68	21.239	3066	3069	3076	rVB	1048655	1350736	1.83%	0.697%
69	21.445	3101	3104	3110	rVB	1035236	1214448	1.65%	0.627%
70	21.533	3115	3119	3128	rVB	2857034	4090761	5.55%	2.111%
71	23.797	3498	3504	3510	rVB2	2316357	4386543	5.95%	2.264%
72	23.950	3525	3530	3539	rVB	1861926	3922991	5.32%	2.024%

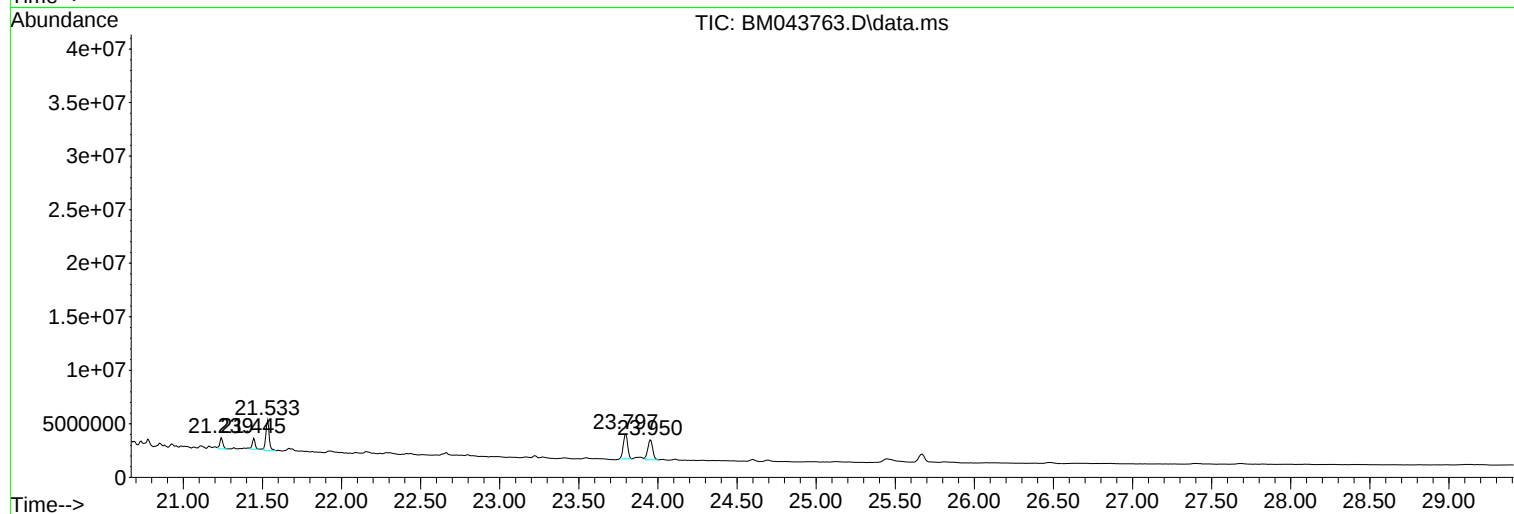
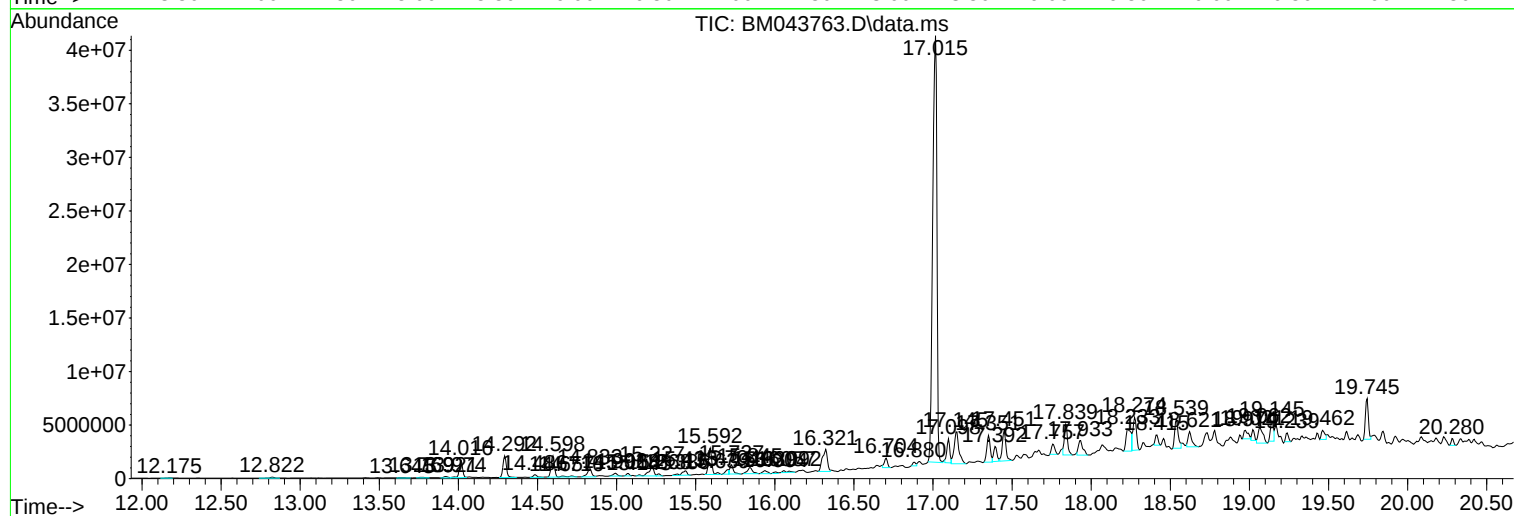
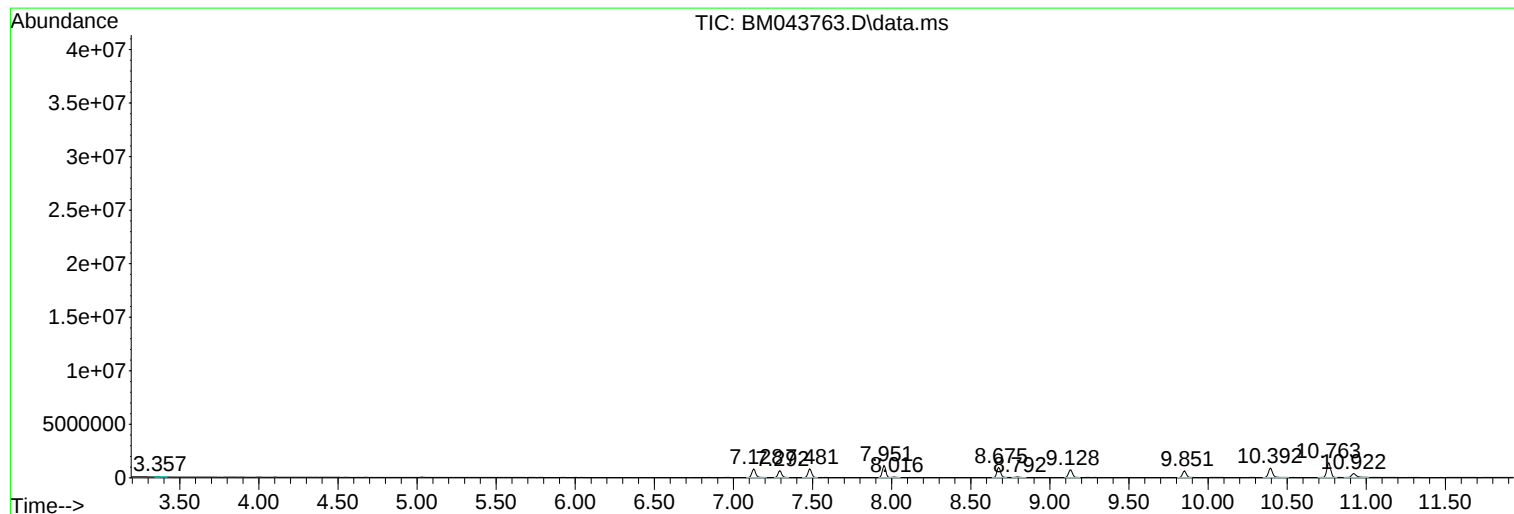
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Instrument :
BNA_M
ClientSampleId :
GCF90ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

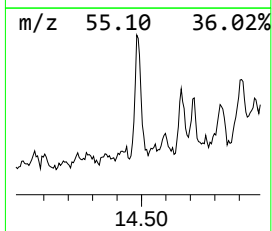
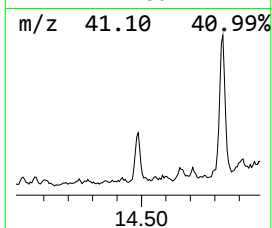
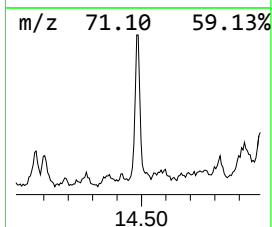
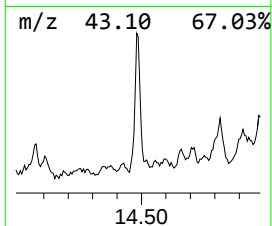
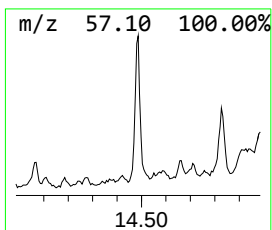
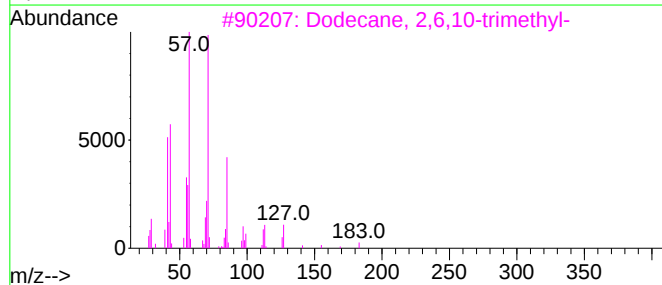
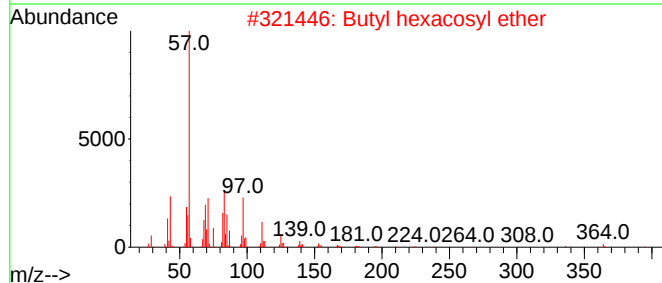
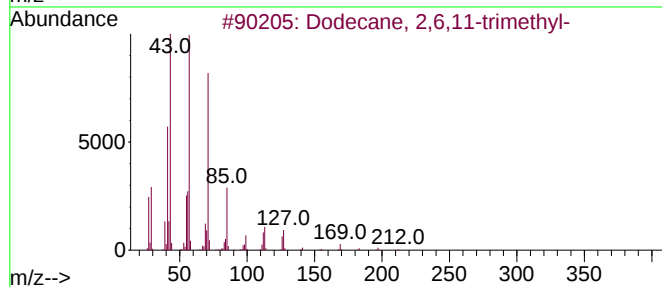
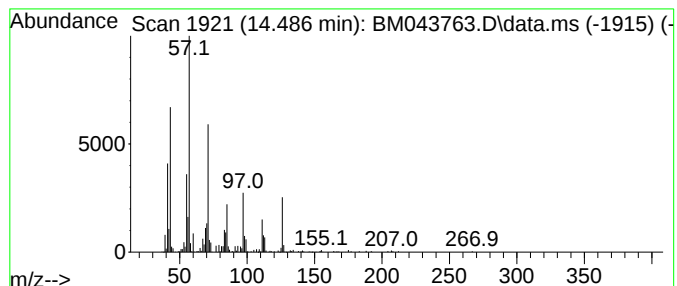
Instrument :
BNA_M
ClientSampleId :
GCF90ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.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 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.486	3.15 ng/ul	462557	Acenaphthene-d10	14.598	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	60
2	Butyl hexacosyl ether	438	C30H62O	1000406-41-0	53
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	52
4	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	50
5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	50



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Instrument :
BNA_M
ClientSampleId :
GCF90ME

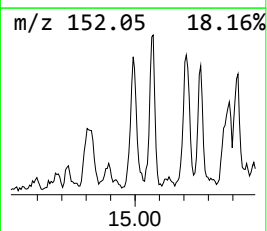
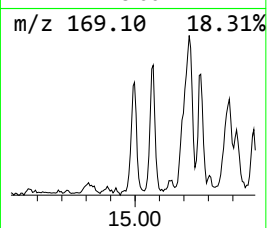
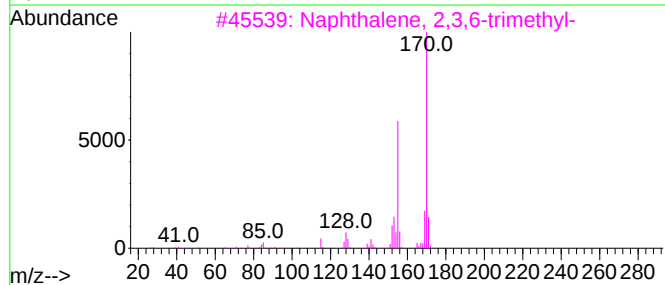
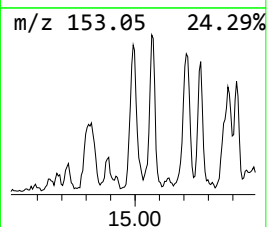
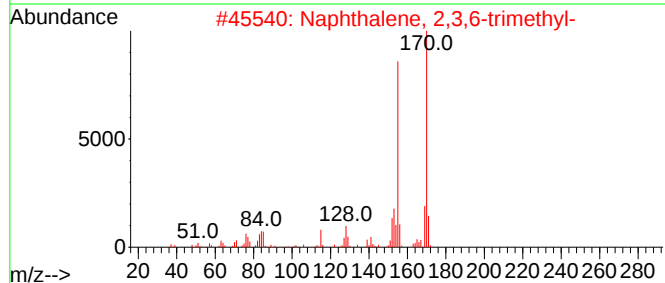
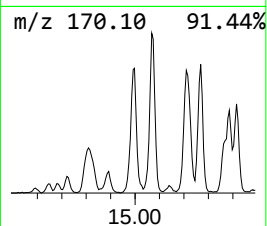
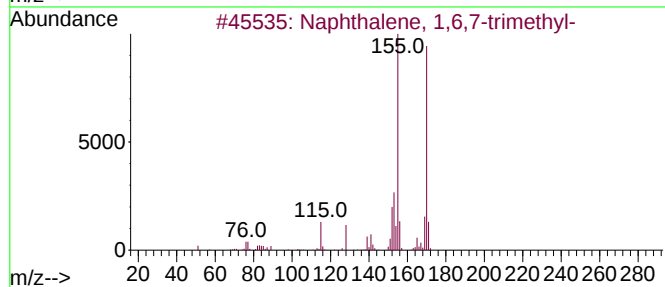
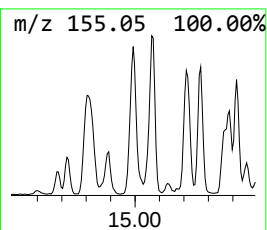
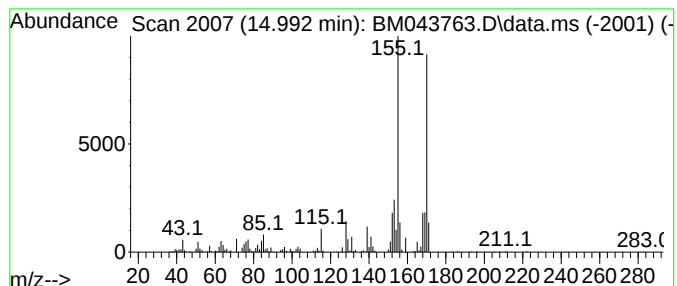
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Naphthalene, 1,6,7-trimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.992	3.14 ng/ul	460439	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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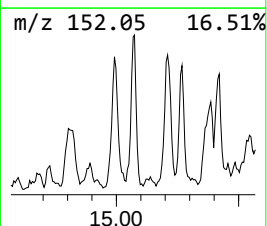
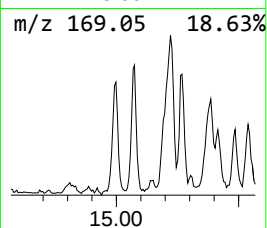
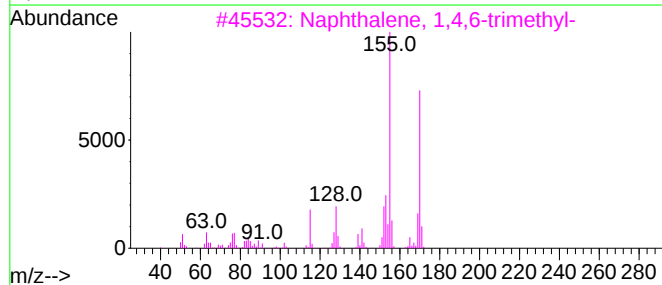
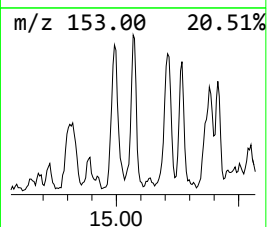
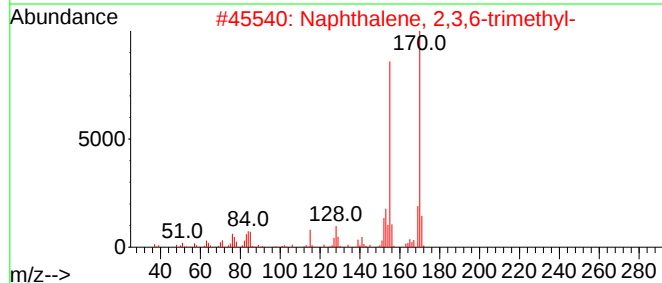
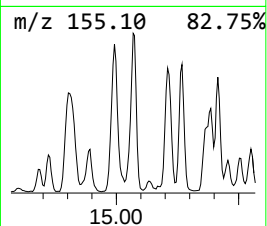
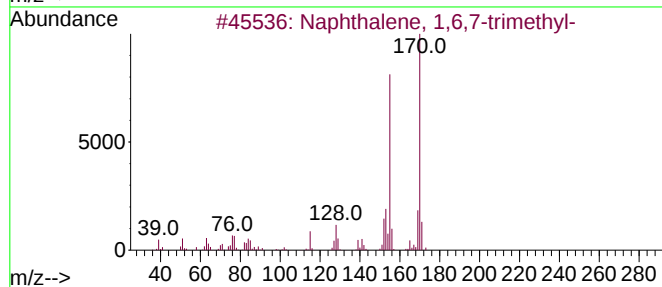
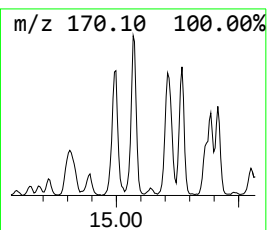
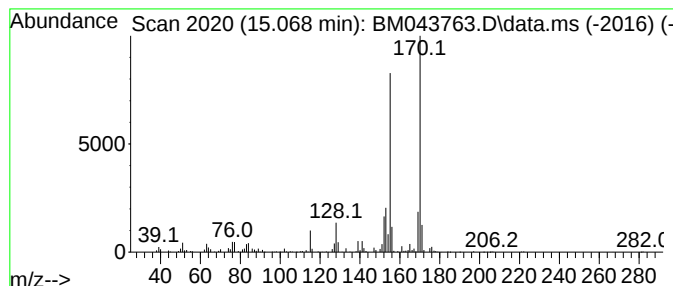
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 2,3,6-trimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.069	2.05 ng/ul	300986	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94



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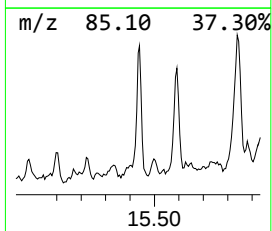
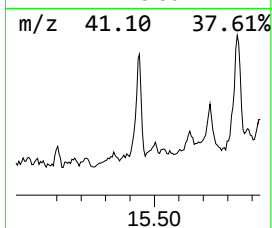
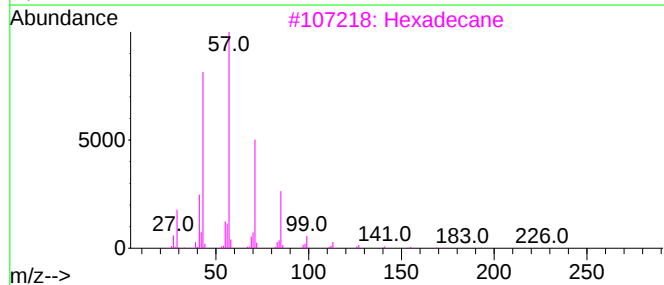
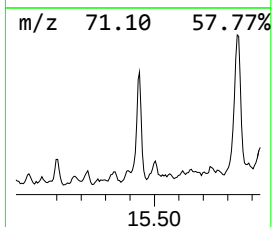
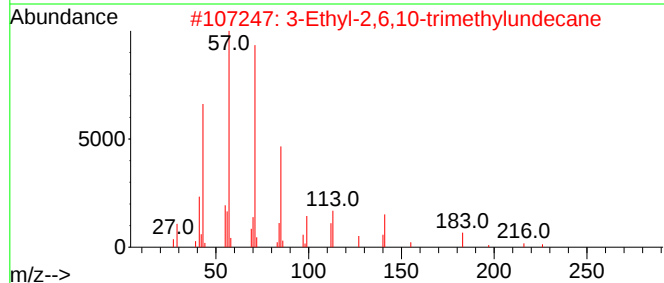
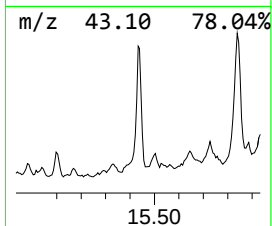
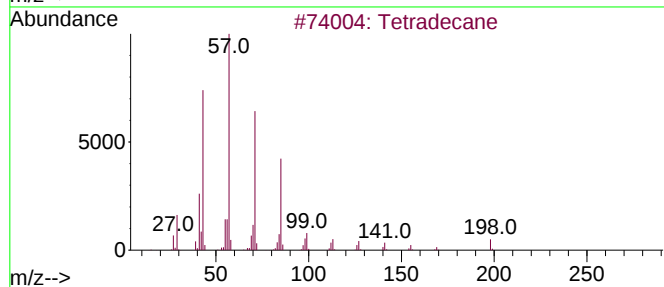
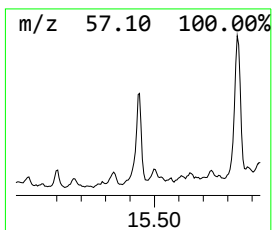
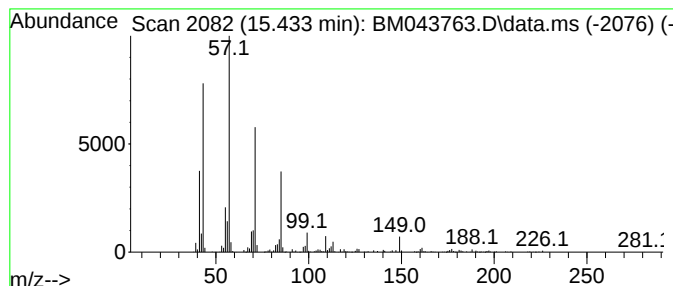
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	5.66 ng/ul	830436	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	90
3			Hexadecane	226	C16H34	000544-76-3	86
4			Heptadecane	240	C17H36	000629-78-7	86
5			Tridecane	184	C13H28	000629-50-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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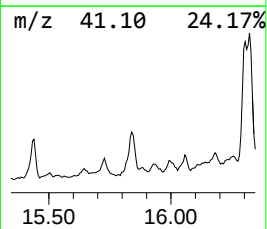
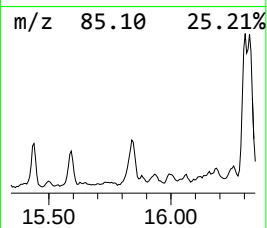
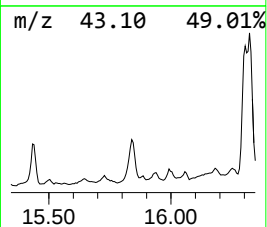
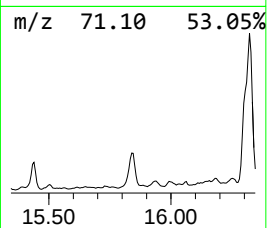
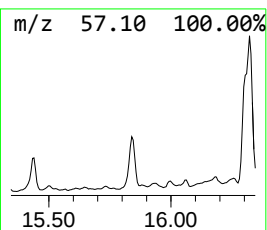
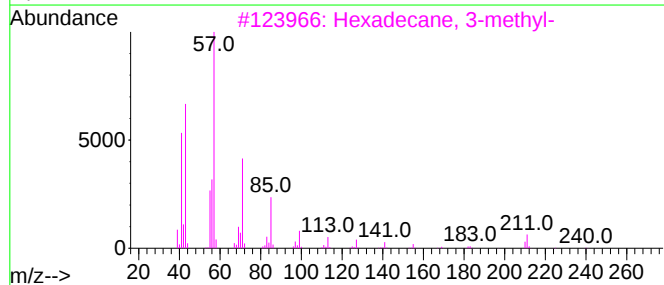
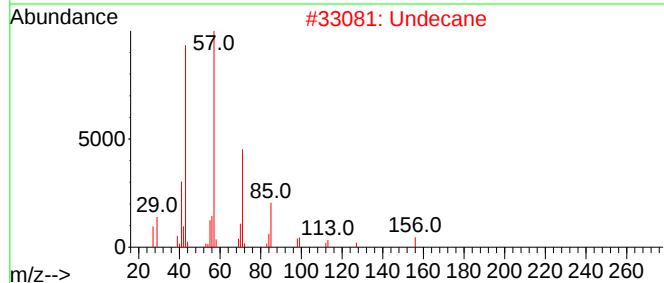
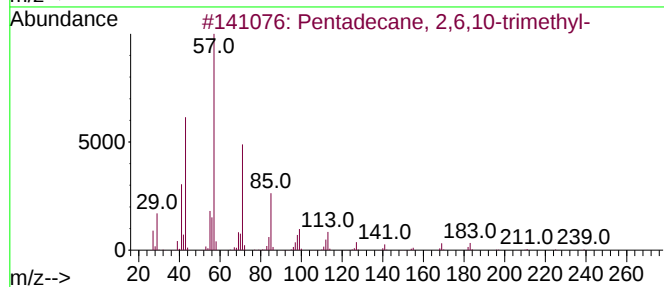
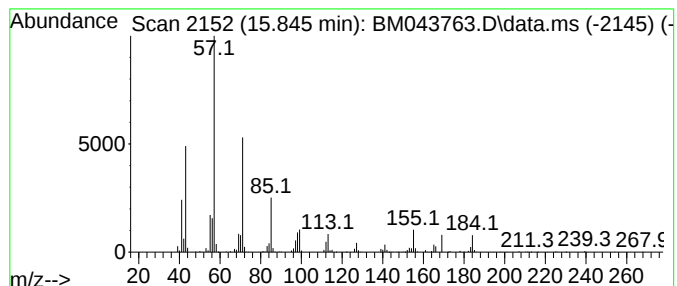
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.845	8.80 ng/ul	1291550	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2			Undecane	156	C11H24	001120-21-4	74
3			Hexadecane, 3-methyl-	240	C17H36	006418-43-5	68
4			Heptacosane	380	C27H56	000593-49-7	64
5			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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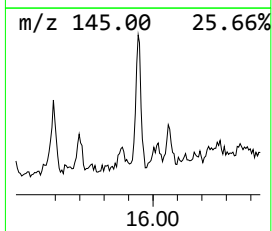
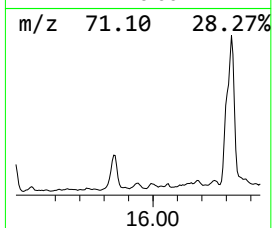
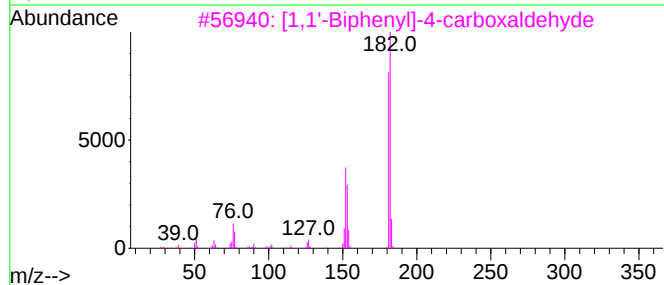
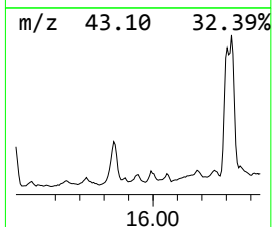
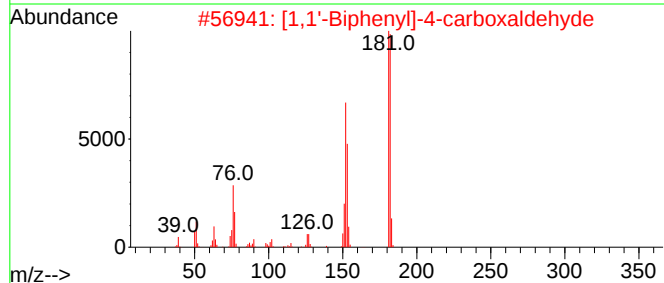
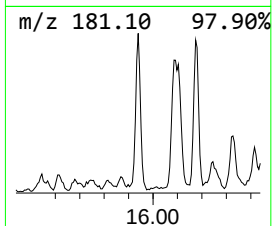
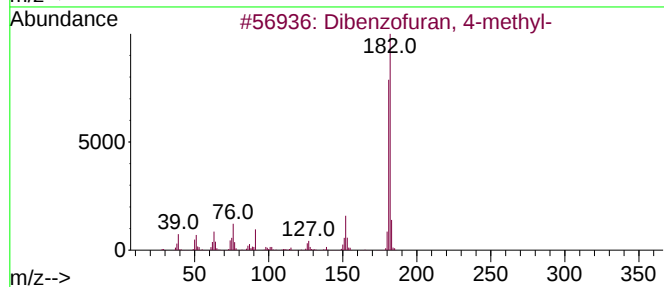
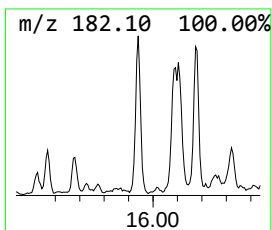
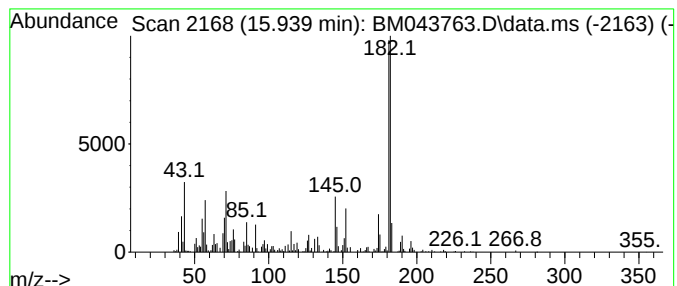
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Dibenzofuran, 4-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.939	3.91 ng/ul	574044	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	89
2			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	59
3			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	53
4			3-(2-Naphthyl)acrylaldehyde	182	C13H10O	020884-05-3	53
5			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
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ClientSampleId :
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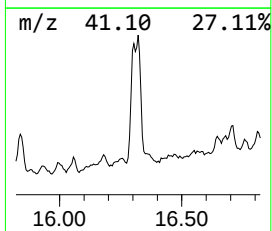
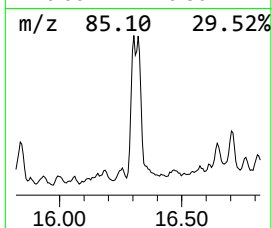
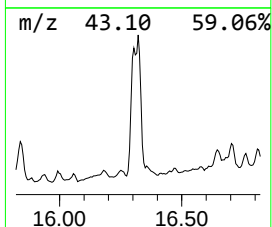
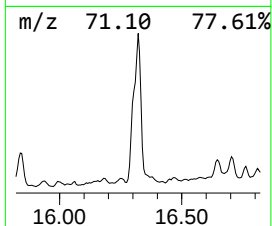
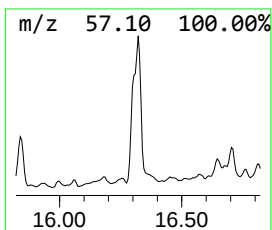
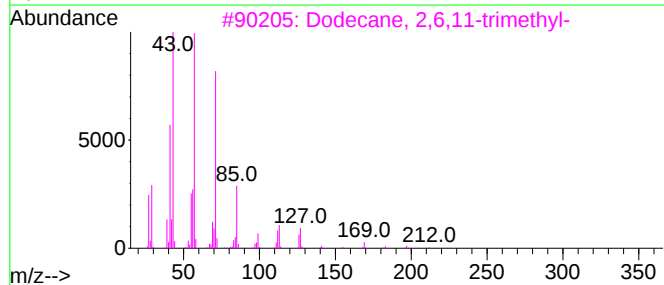
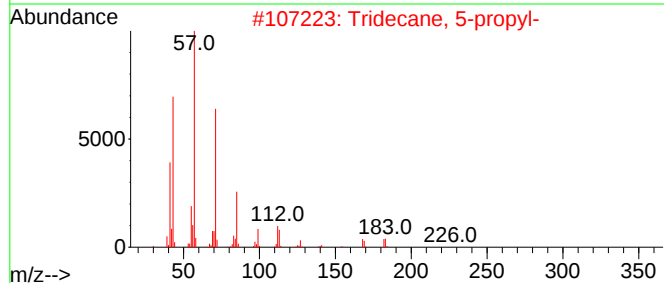
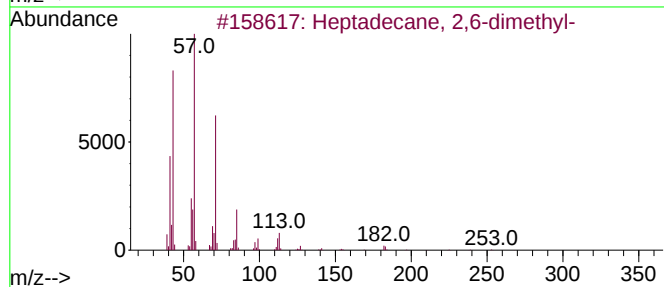
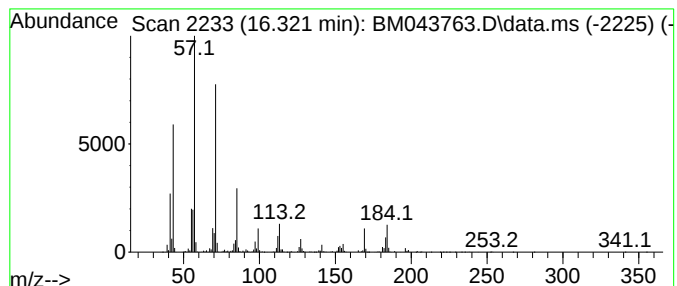
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.321	23.23 ng/ul	4474630	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	91
2		Tridecane, 5-propyl-	226	C16H34	055045-11-9	81
3		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5		Tridecane, 7-methyl-	198	C14H30	026730-14-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
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Misc :
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Instrument :
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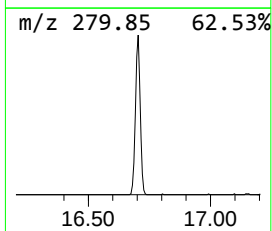
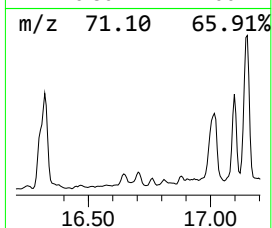
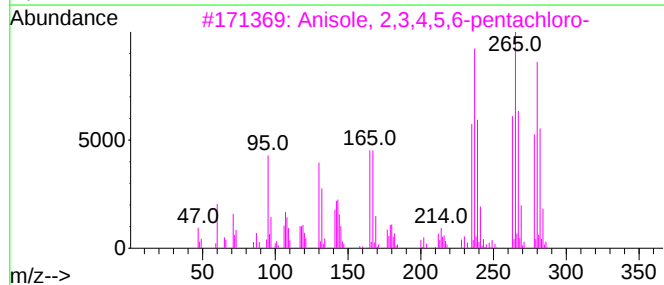
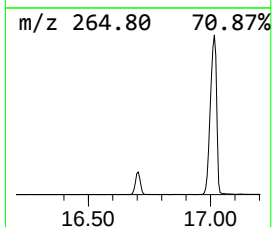
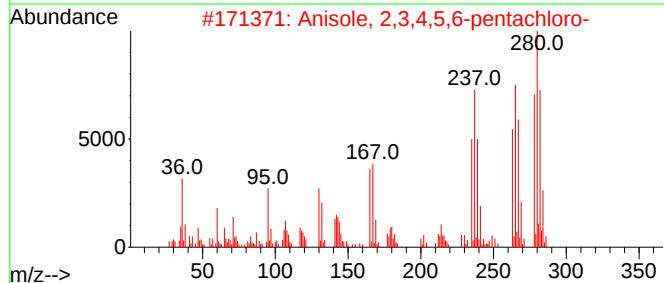
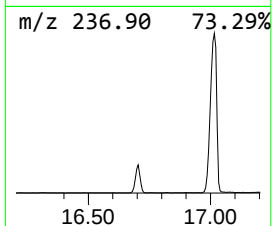
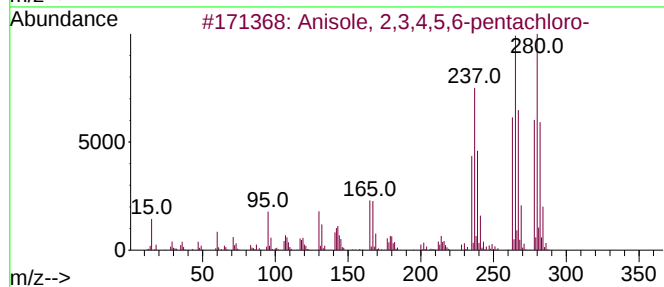
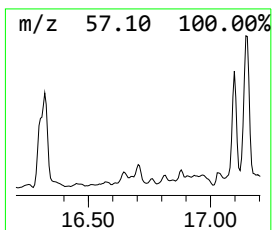
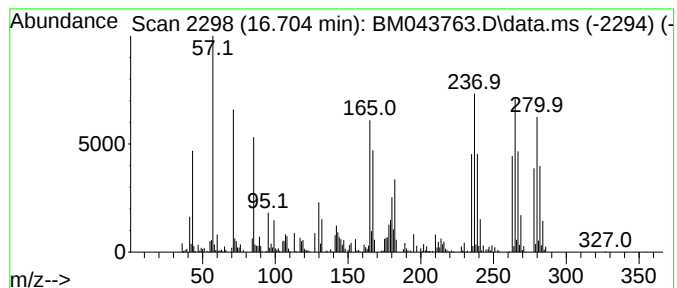
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Anisole, 2,3,4,5,6-pentachl... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.704	6.25 ng/ul	1204830	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	99		
2	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	94		
3	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	91		
4	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	64		
5	3,5-Dibromobenzoic acid	278	C7H4Br2O2	000618-58-6	42		



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Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
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Sample : 05952-07ME
Misc :
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Instrument :
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ClientSampleId :
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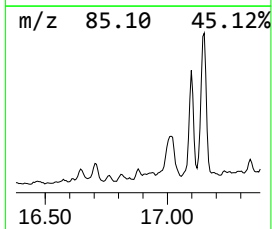
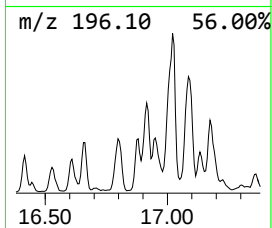
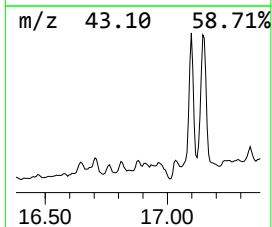
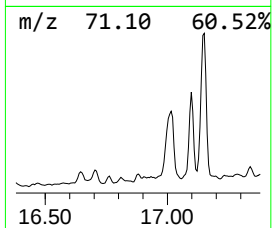
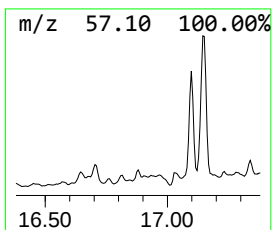
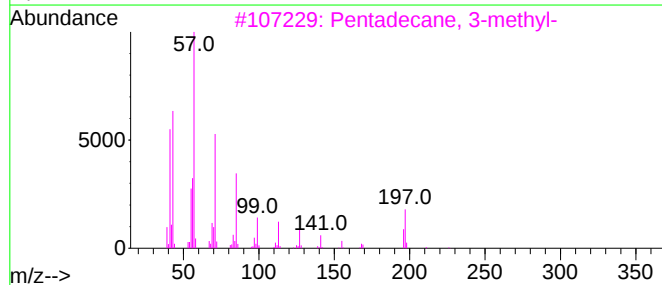
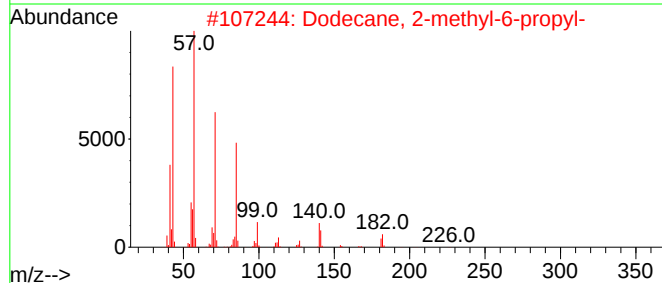
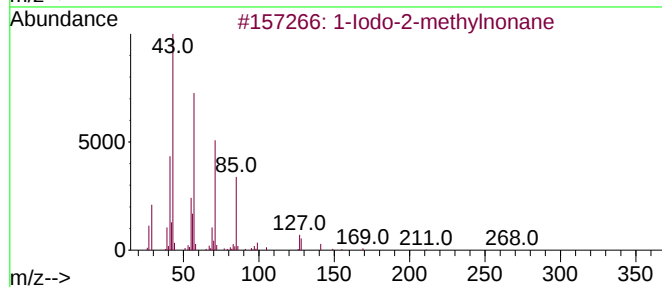
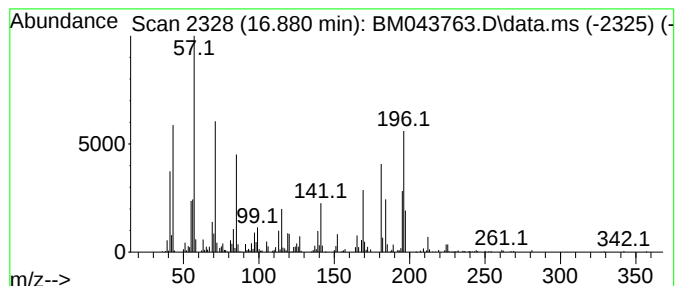
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.880	2.54 ng/ul	489227	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	35
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	30
3			Pentadecane, 3-methyl-	226	C16H34	002882-96-4	25
4			Heptadecane, 3-methyl-	254	C18H38	006418-44-6	25
5			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Misc :
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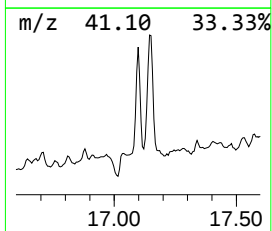
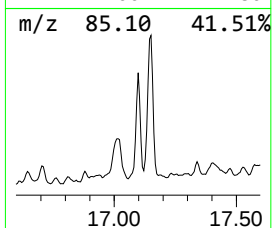
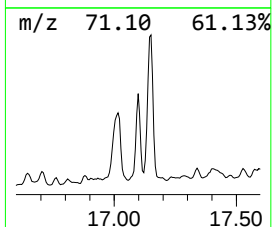
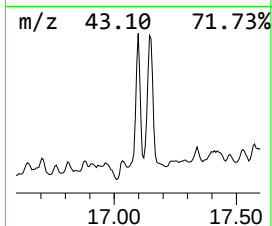
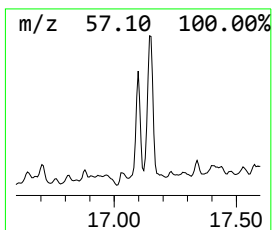
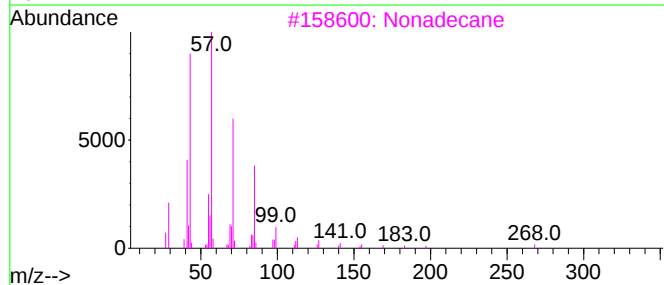
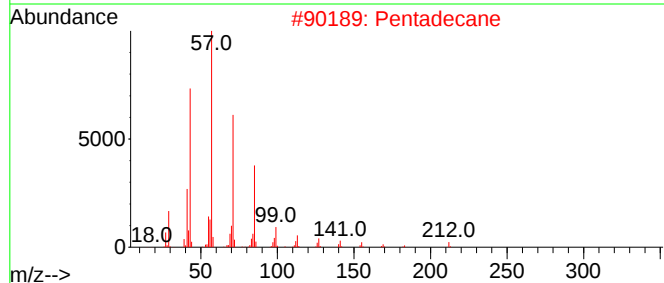
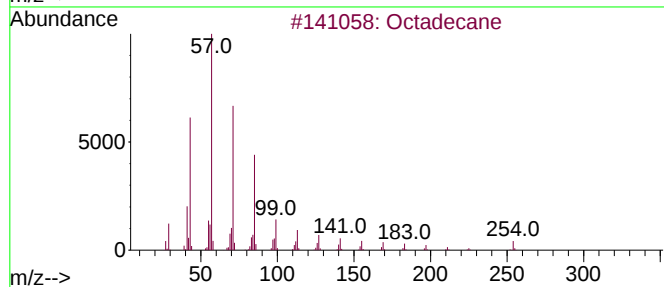
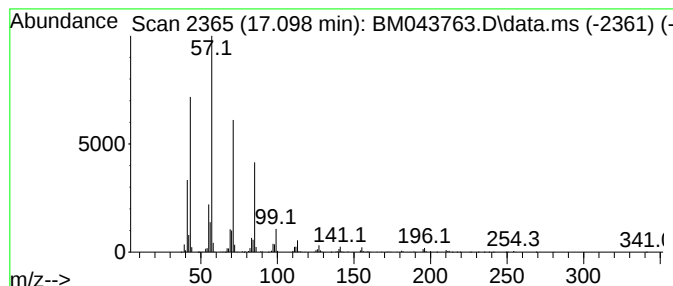
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.098	14.91 ng/ul	2871720	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	95
2		Pentadecane	212	C15H32	000629-62-9	91
3		Nonadecane	268	C19H40	000629-92-5	91
4		Heptacosane	380	C27H56	000593-49-7	90
5		Heptadecane	240	C17H36	000629-78-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
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Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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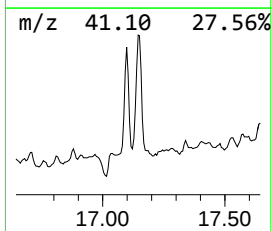
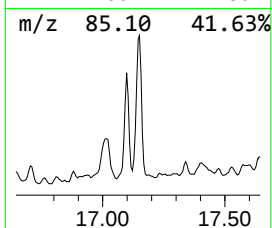
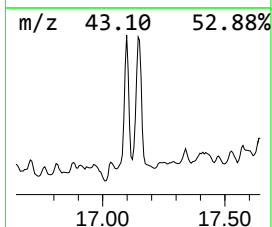
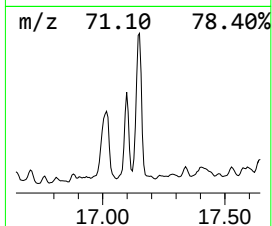
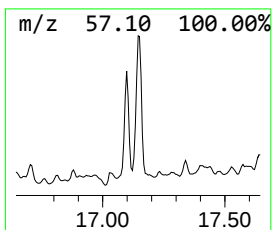
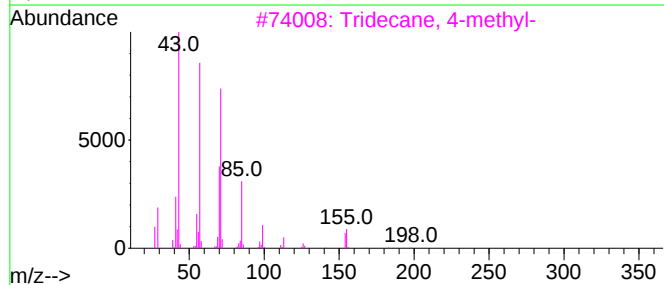
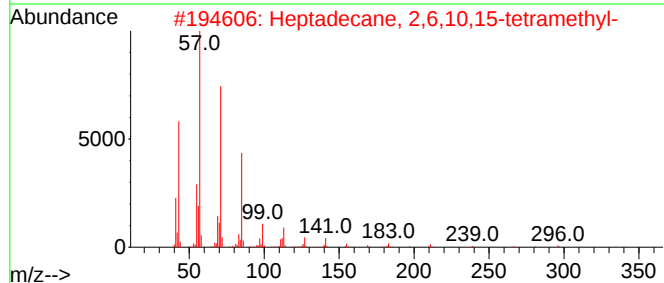
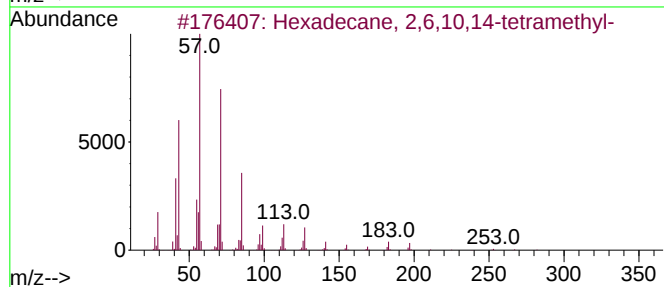
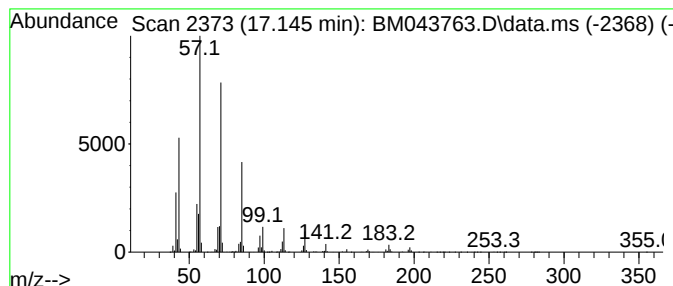
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.145	32.40 ng/ul	6242350	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
3			Tridecane, 4-methyl-	198	C14H30	026730-12-1	87
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5			Dodecane	170	C12H26	000112-40-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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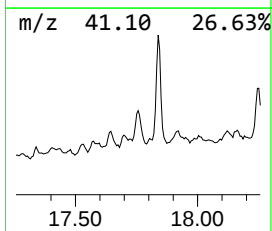
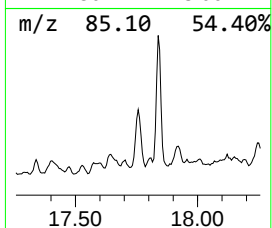
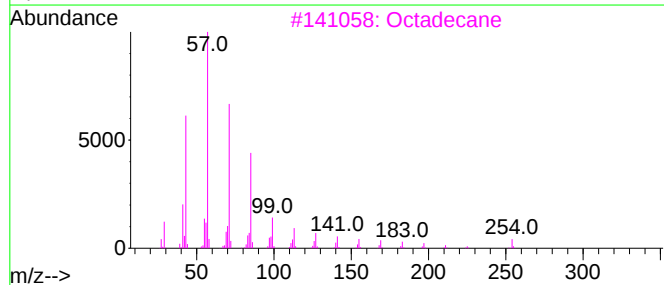
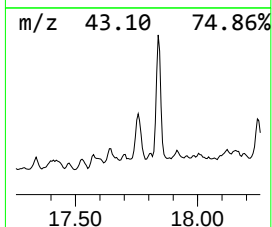
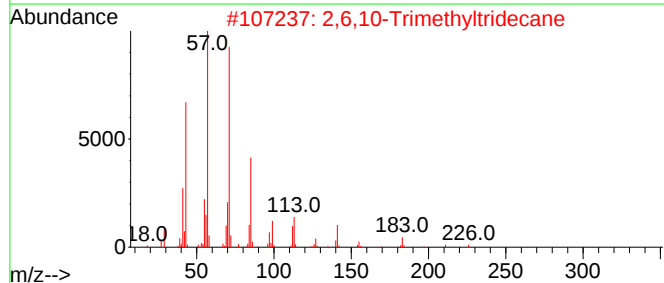
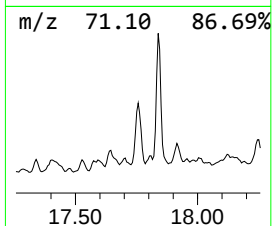
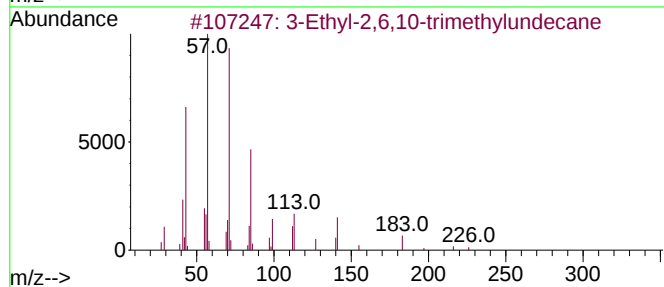
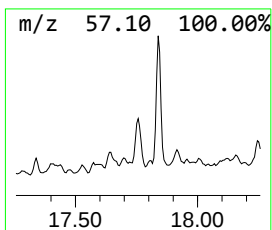
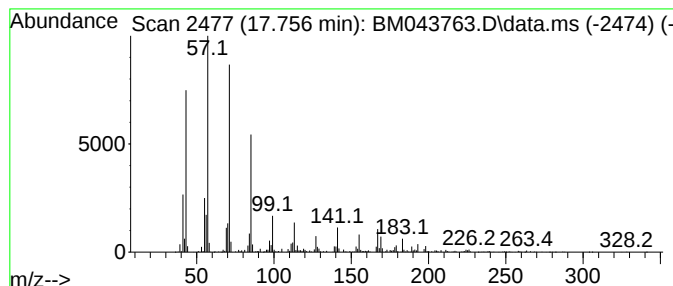
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.756	7.32 ng/ul	1410860	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	89
2		2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	89
3		Octadecane	254	C18H38	000593-45-3	87
4		Hentriacontane	437	C31H64	000630-04-6	86
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
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Instrument :
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ClientSampleId :
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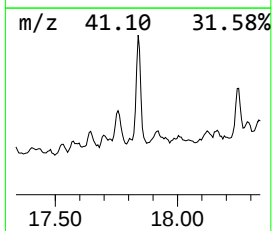
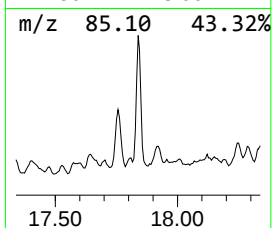
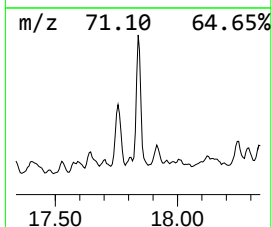
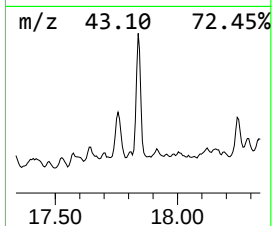
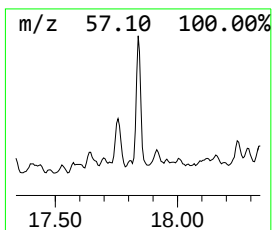
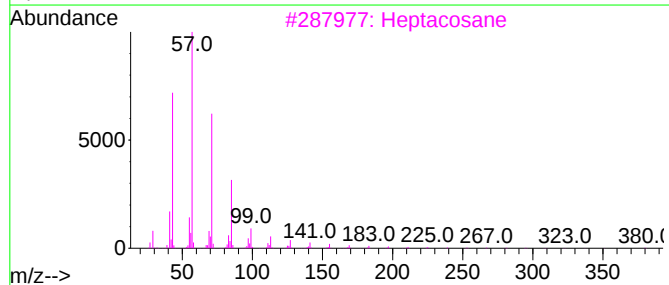
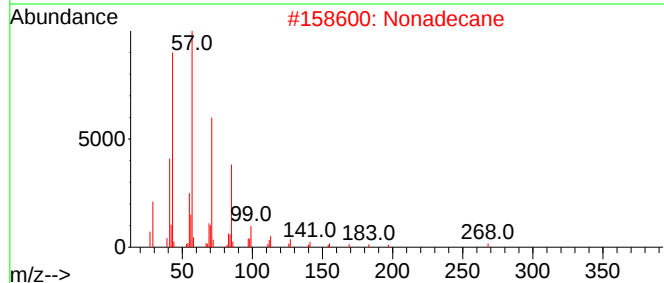
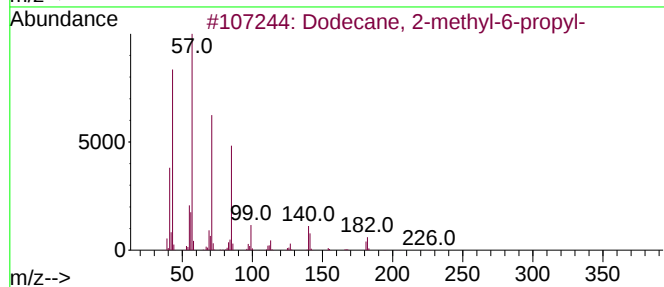
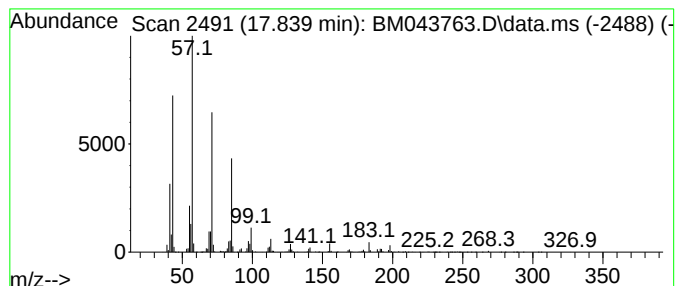
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.839	20.59 ng/ul	3967670	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91
2		Nonadecane	268	C19H40	000629-92-5	90
3		Heptacosane	380	C27H56	000593-49-7	90
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5		Pentadecane	212	C15H32	000629-62-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Acq On : 05 Jan 2024 16:49
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ClientSampleId :
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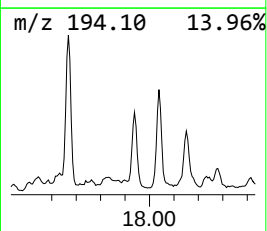
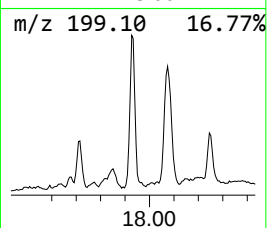
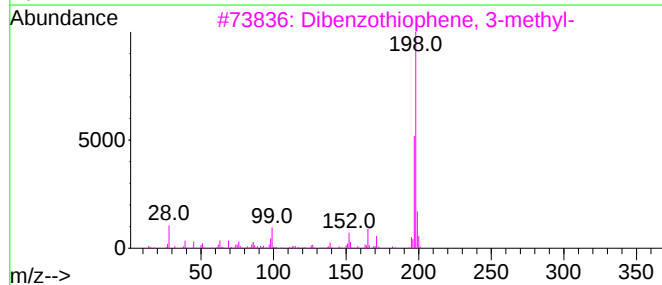
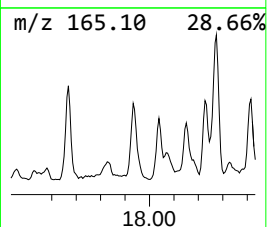
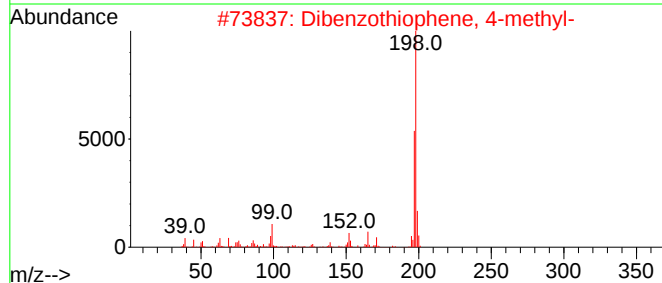
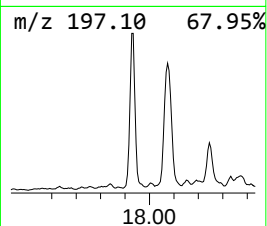
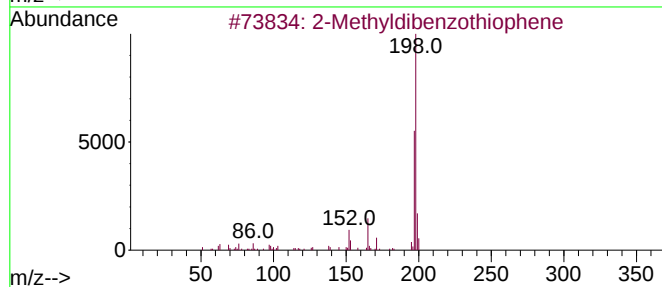
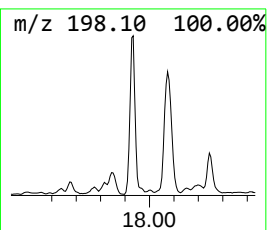
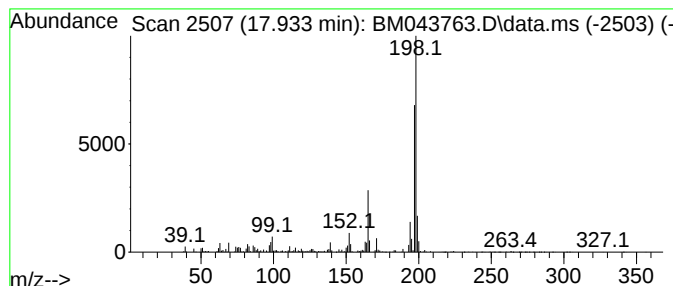
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 2-Methyldibenzothiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.933	13.47 ng/ul	2595650	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	90
2			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	81
3			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	81
4			Thioxanthene	198	C13H10S	000261-31-4	64
5			Tacrine	198	C13H14N2	000321-64-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Instrument :
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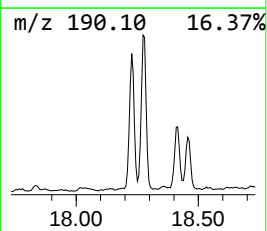
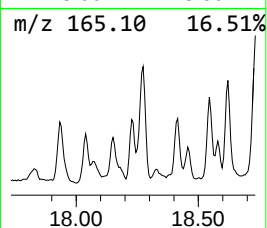
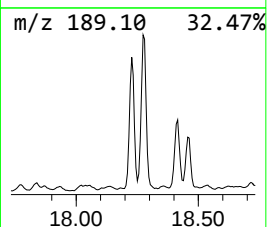
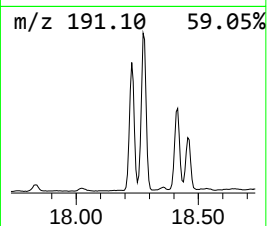
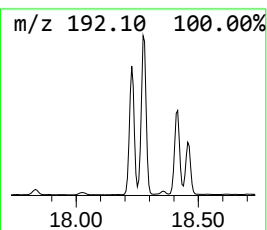
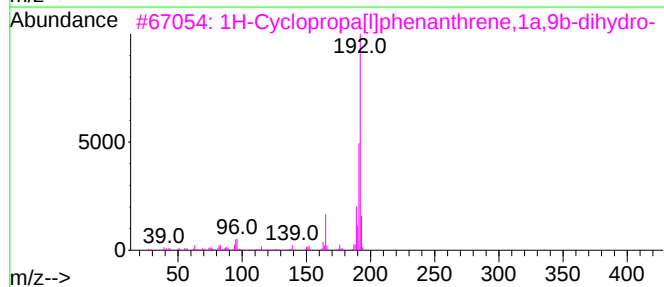
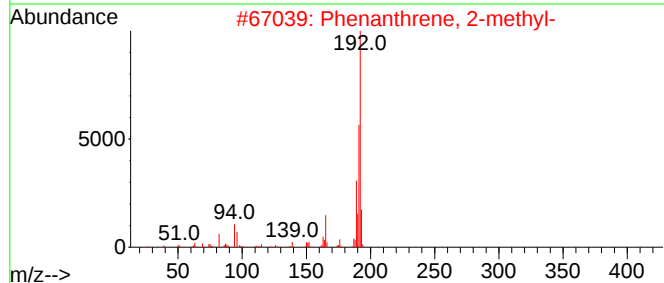
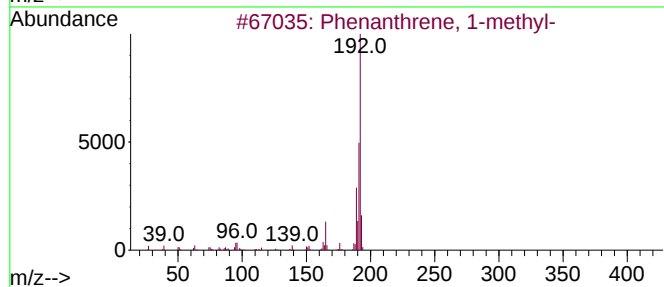
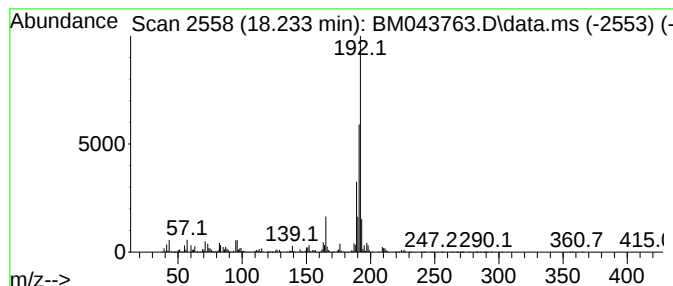
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Phenanthrene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	24.76 ng/ul	4770310	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Acq On : 05 Jan 2024 16:49
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Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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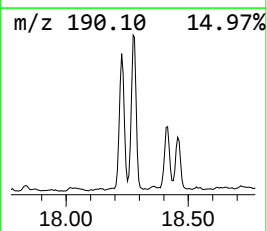
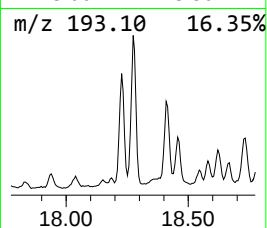
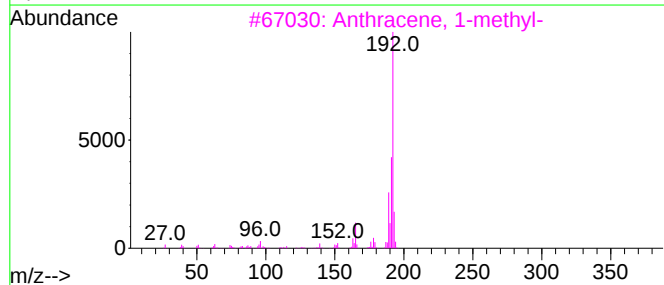
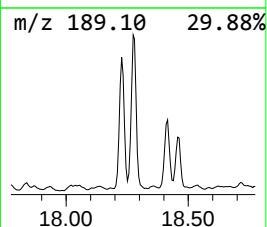
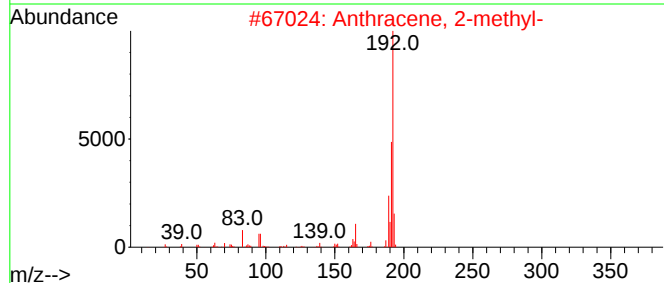
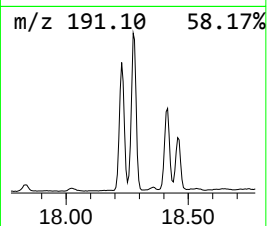
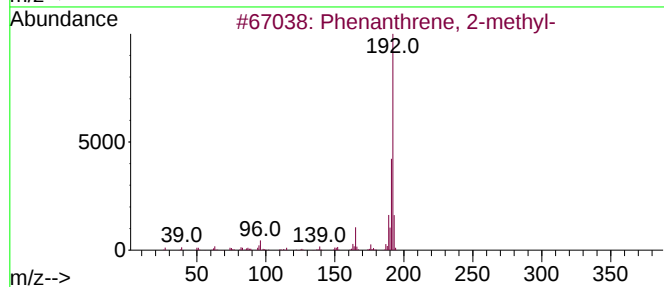
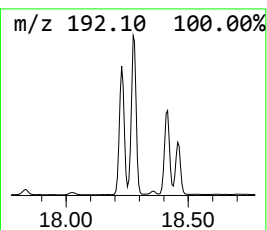
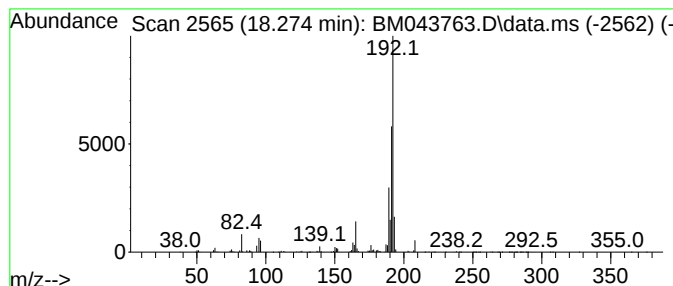
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.274	23.53 ng/ul	4533000	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
4		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Misc :
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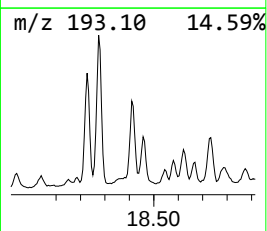
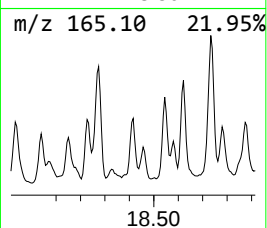
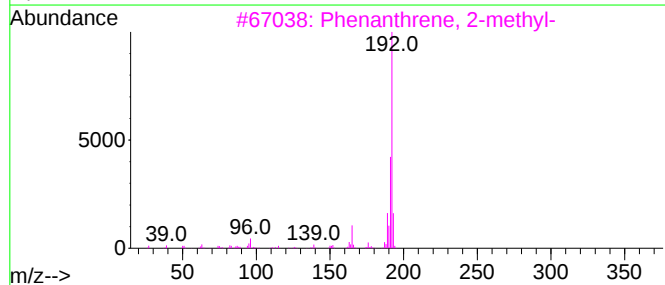
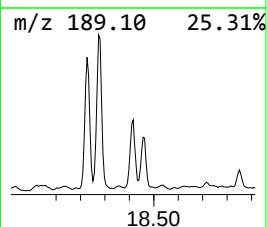
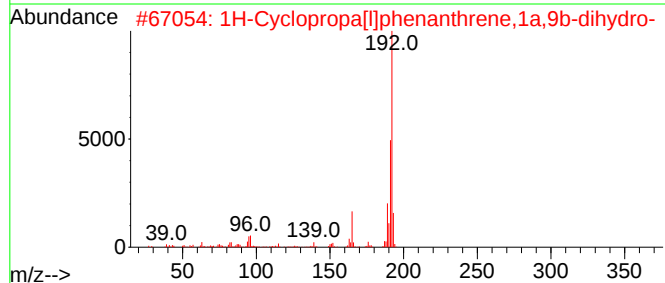
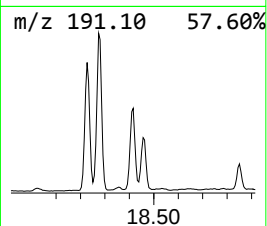
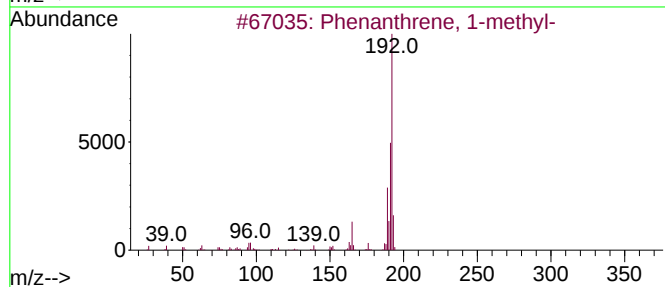
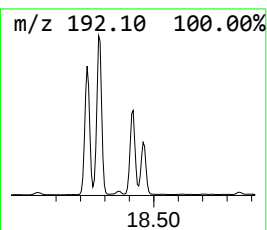
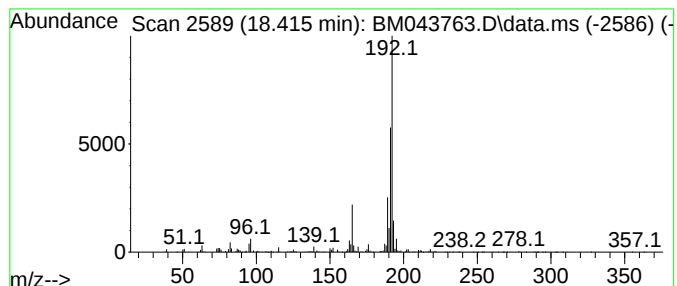
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1H-Cyclopropa[l]phenanthren... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.415	6.16 ng/ul	1186420	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	93
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
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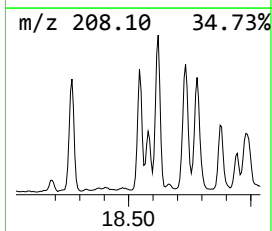
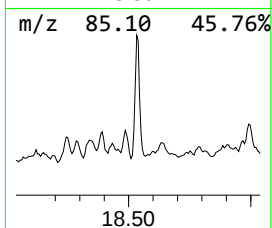
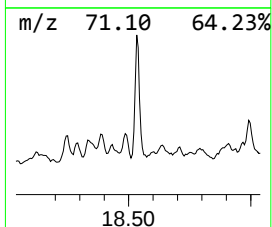
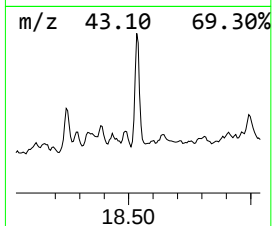
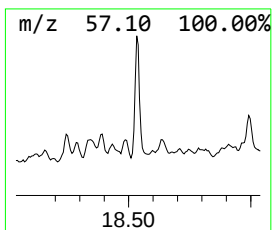
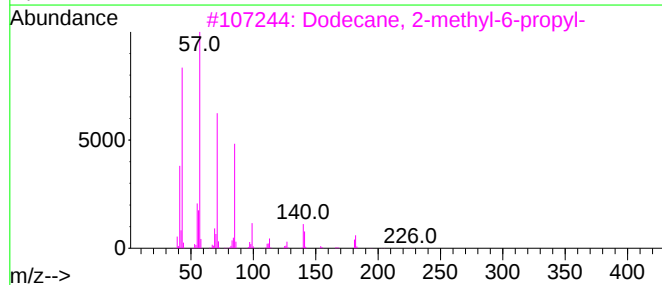
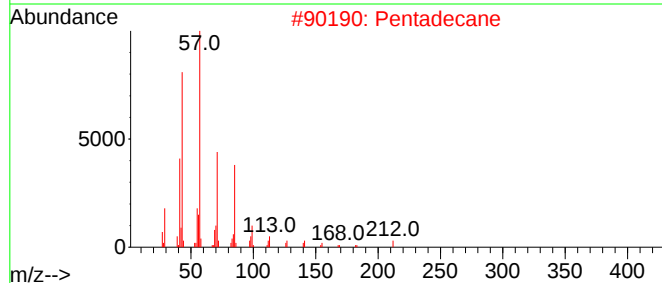
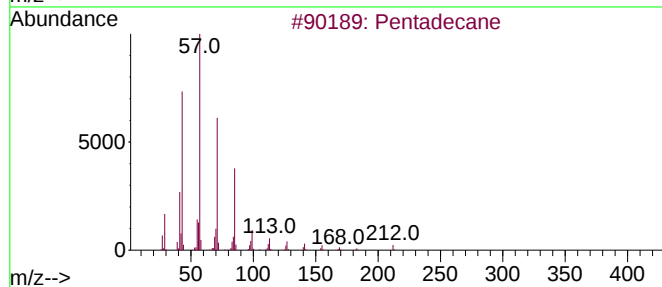
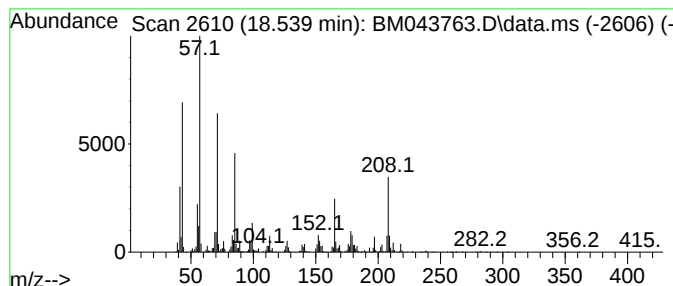
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.539	21.79 ng/ul	4198870	Phenanthrene-d10		17.351	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	97
2	Pentadecane		212	C15H32	000629-62-9	95
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	78
4	Heneicosane		296	C21H44	000629-94-7	60
5	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Sample : 05952-07ME
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Instrument :
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ClientSampleId :
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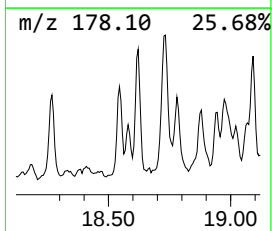
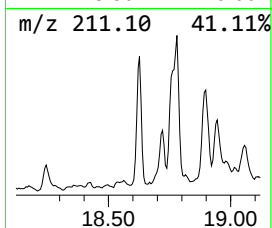
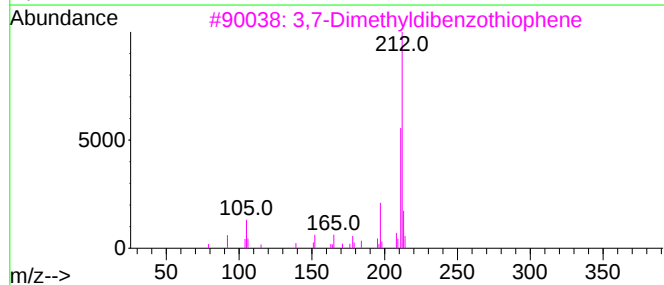
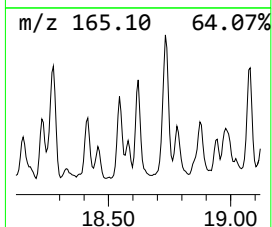
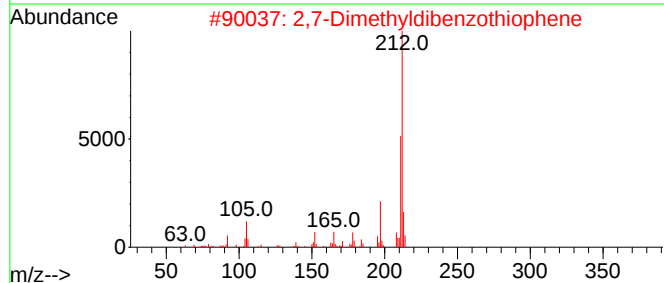
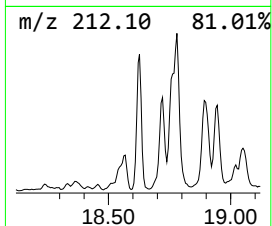
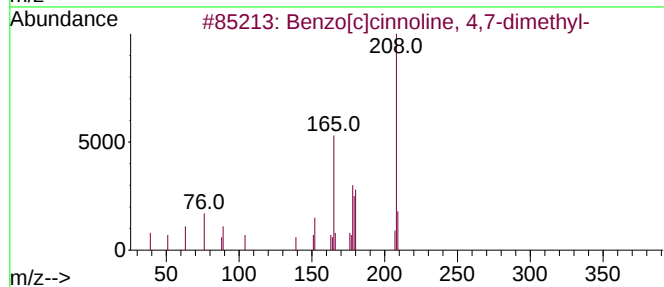
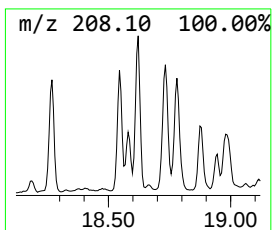
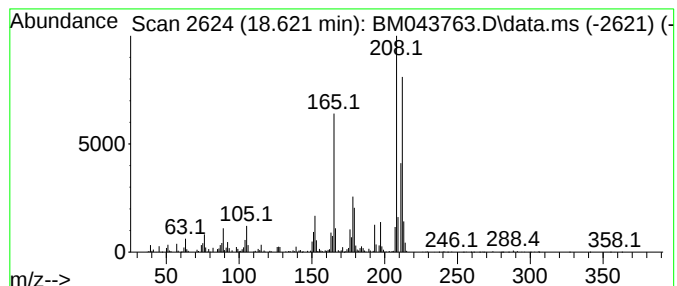
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzo[c]cinnoline, 4,7-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.621	11.99 ng/ul	2309940	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]cinnoline, 4,7-dimethyl-	208	C14H12N2	020684-54-2	89
2			2,7-Dimethyldibenzothiophene	212	C14H12S	031317-19-8	83
3			3,7-Dimethyldibenzothiophene	212	C14H12S	001136-85-2	83
4			9-Fluorenone, 2.4-dimethyl-	208	C15H12O	002928-39-4	70
5			Pinosylvin	212	C14H12O2	022139-77-1	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
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Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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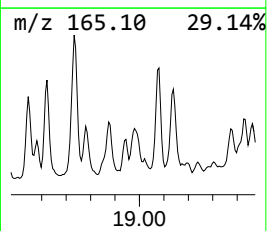
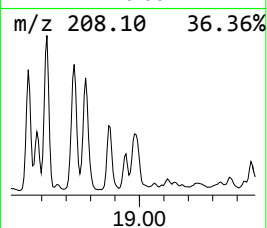
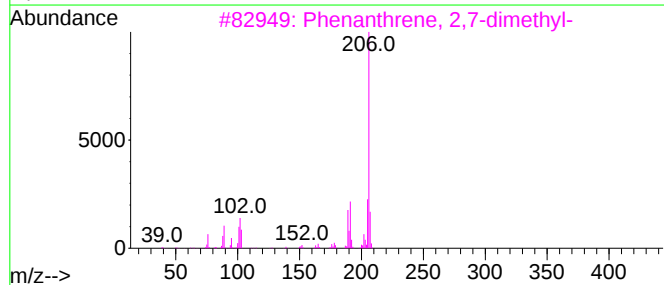
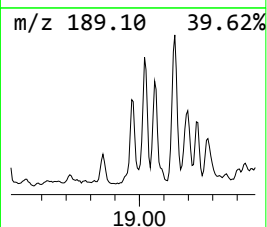
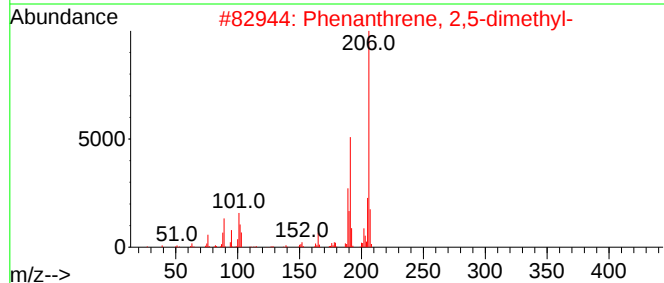
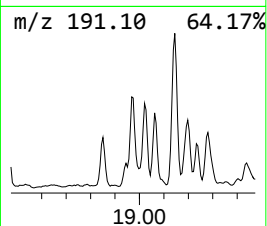
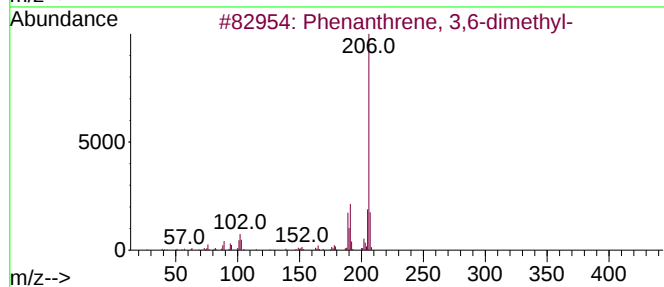
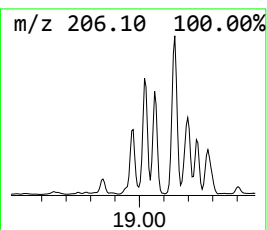
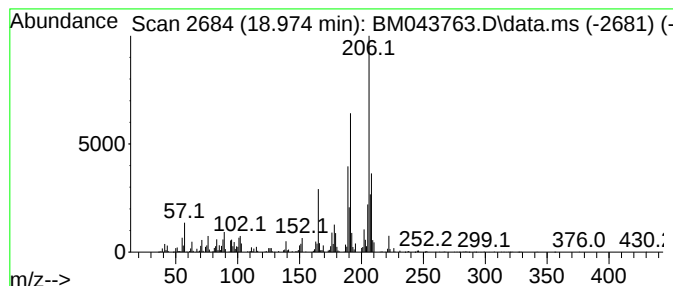
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Phenanthrene, 3,6-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.974	7.92 ng/ul	1525540	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	76
2			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	72
3			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	64
4			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	62
5			Phenanthrene, 9-ethyl-	206	C16H14	003674-75-7	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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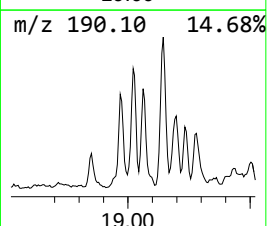
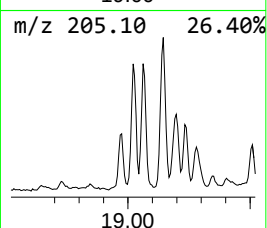
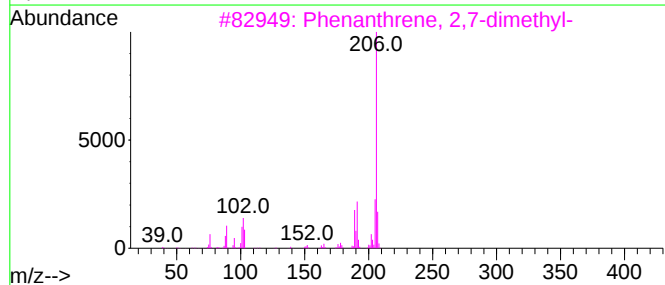
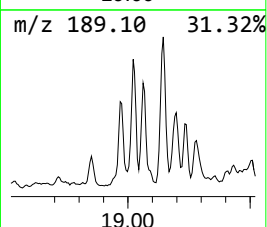
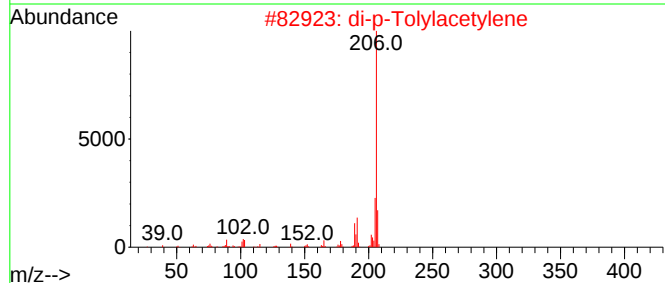
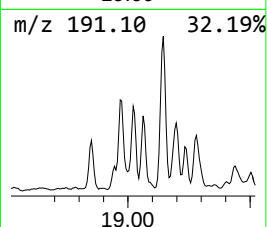
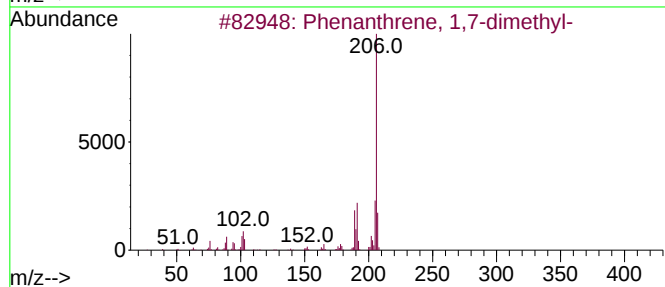
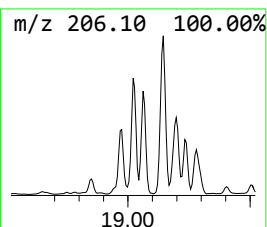
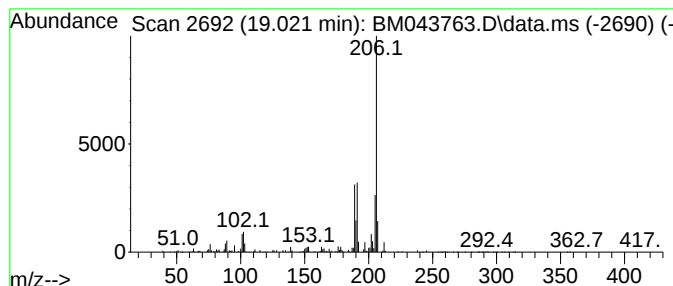
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Phenanthrene, 1,7-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.021	5.41 ng/ul	1041790	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	96
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	94
3			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	93
4			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
5			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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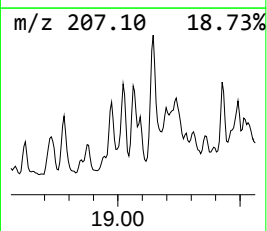
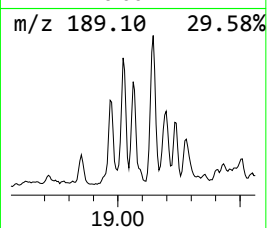
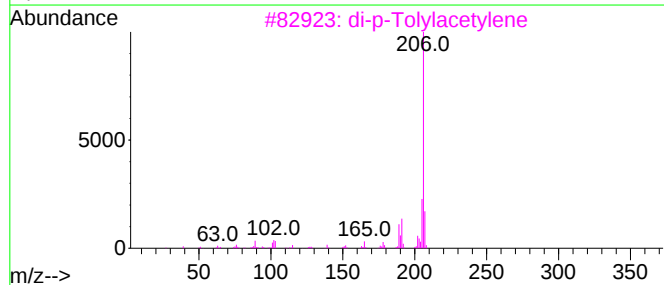
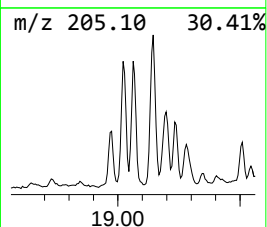
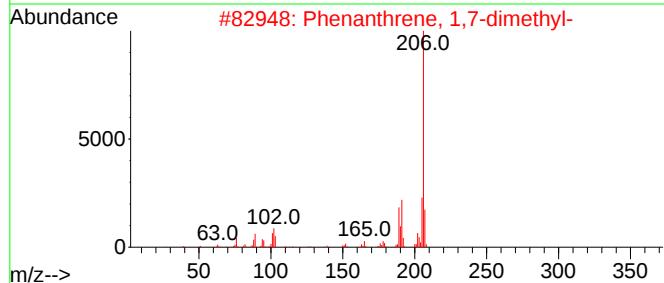
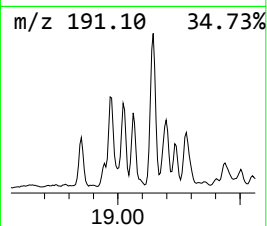
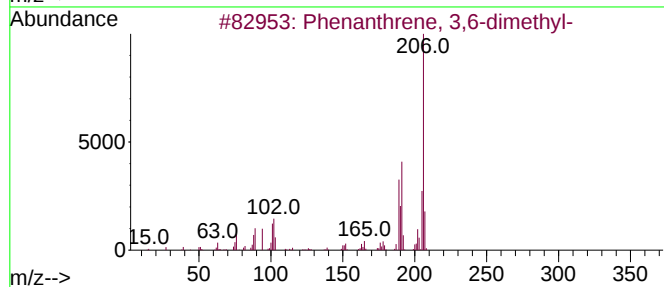
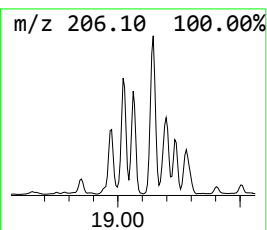
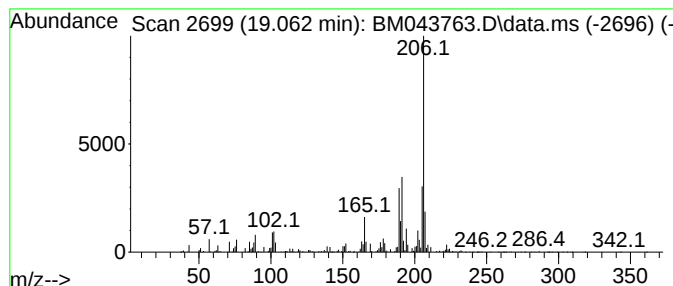
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 di-p-Tolylacetylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.062	16.12 ng/ul	3104910	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	95
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	91
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

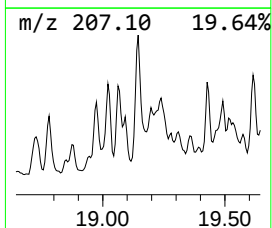
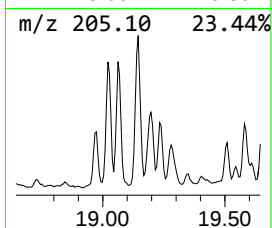
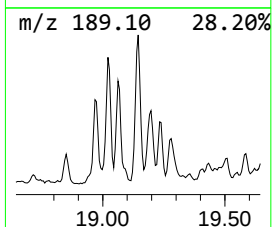
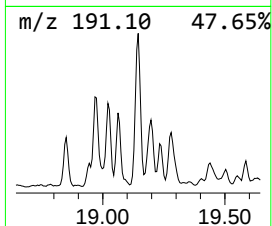
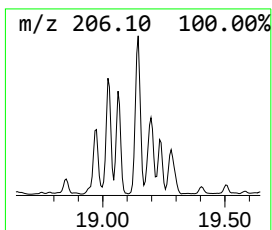
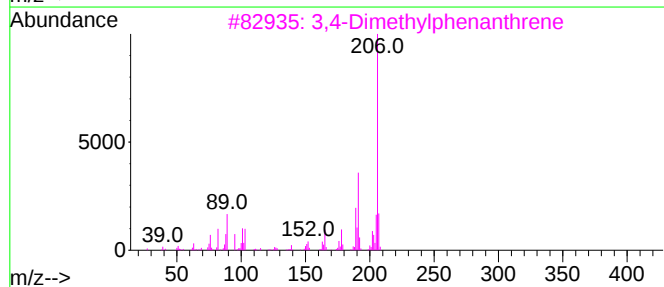
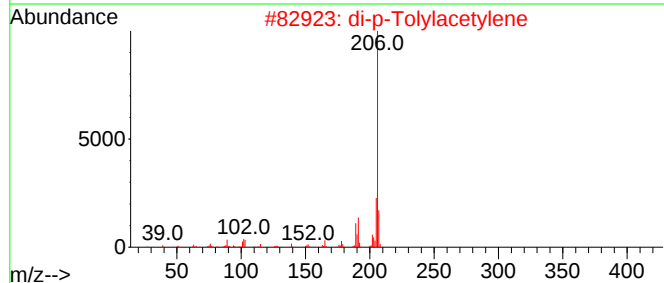
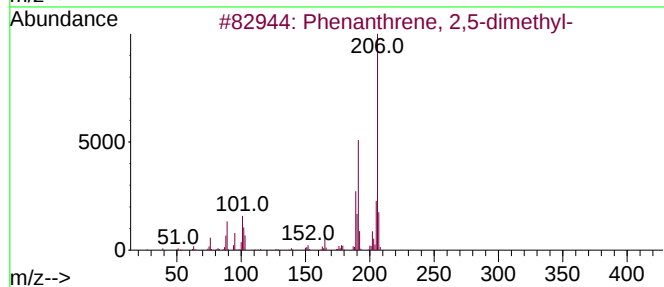
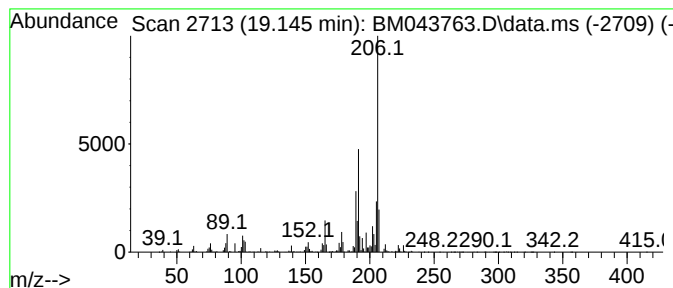
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.145	13.72 ng/ul	2643920	Phenanthrene-d10	17.351	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	93
4	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	89
5	9,10-Dimethylanthracene	206	C16H14	000781-43-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
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Sample : 05952-07ME
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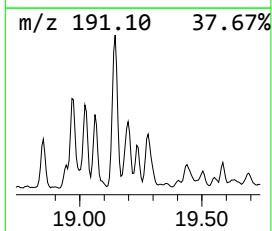
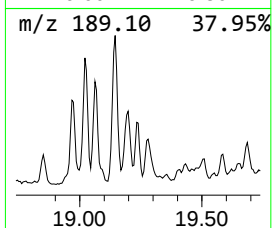
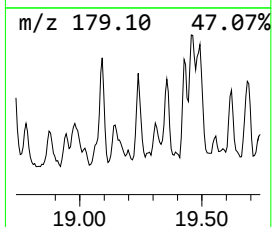
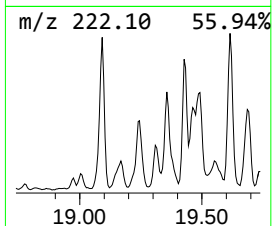
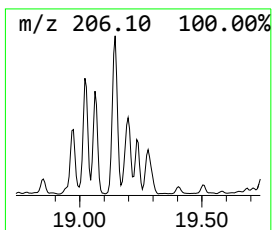
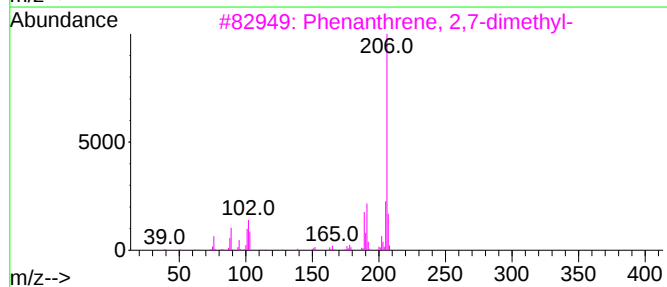
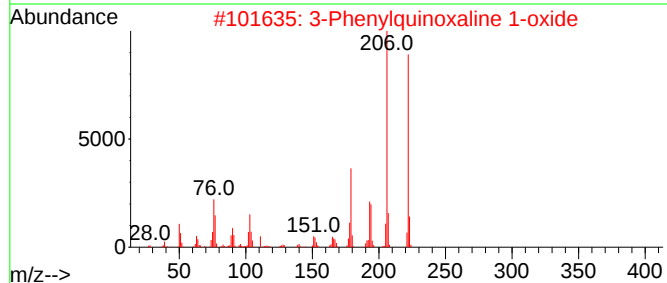
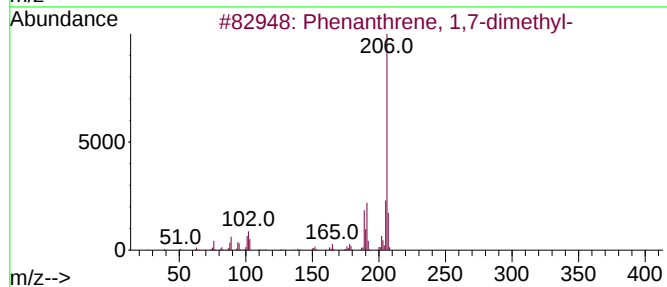
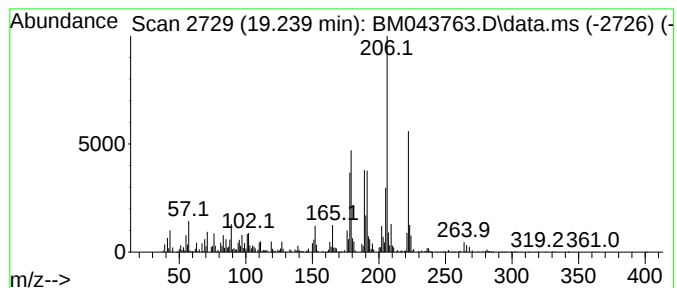
Instrument :
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ClientSampleId :
GCF90ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 3-Phenylquinoxaline 1-oxide Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.239	5.24 ng/ul	1010200	Phenanthrene-d10	17.351	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	64
2	3-Phenylquinoxaline 1-oxide	222	C14H10N2O	010129-97-2	62
3	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	50
4	Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	49
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

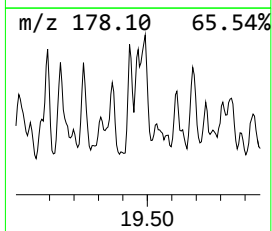
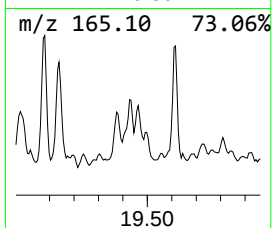
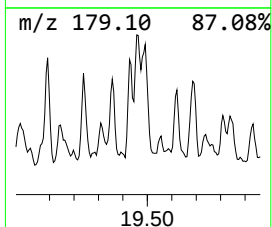
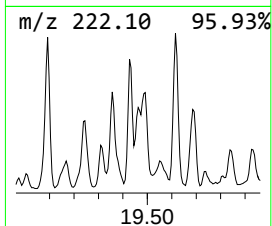
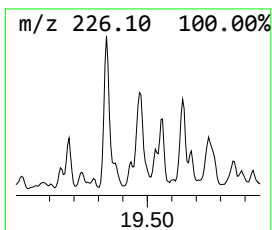
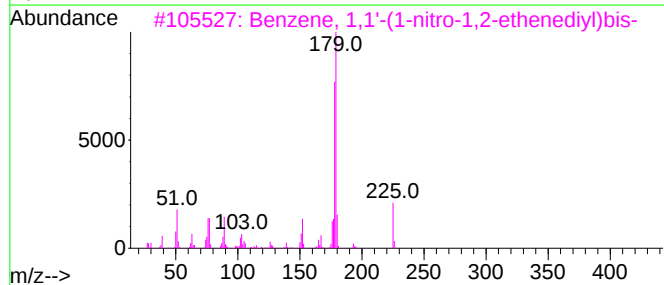
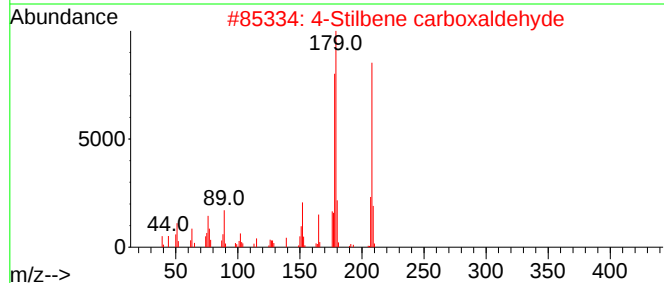
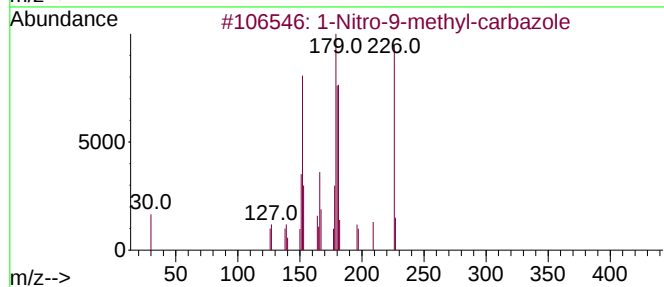
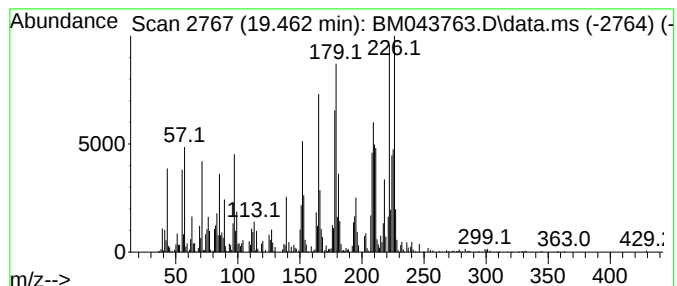
Instrument :
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ClientSampleId :
GCF90ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 1-Nitro-9-methyl-carbazole Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.462	6.87 ng/ul	1406120	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nitro-9-methyl-carbazole	226	C13H10N2O2	066243-81-0	55
2	4-Stilbene carboxaldehyde	208	C15H12O	032555-96-7	35
3	Benzene, 1,1'-(1-nitro-1,2-ethen...	225	C14H11NO2	001215-07-2	35
4	3,4-diphenyl-3-buten-2-one	222	C16H14O	1000378-97-9	30
5	1-Propene-2-thiol, 1,1-diphenyl-	226	C15H14S	074630-83-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

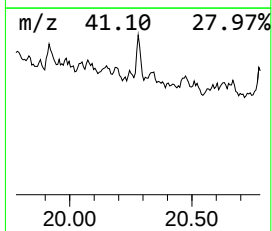
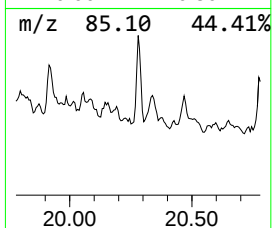
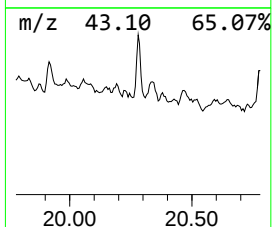
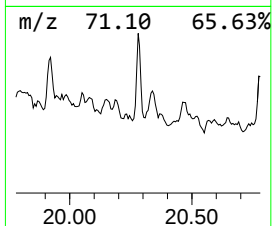
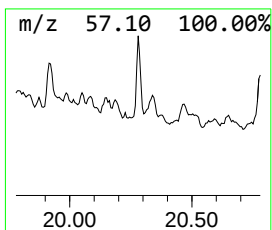
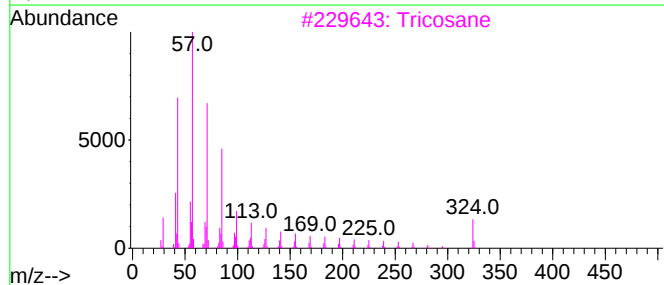
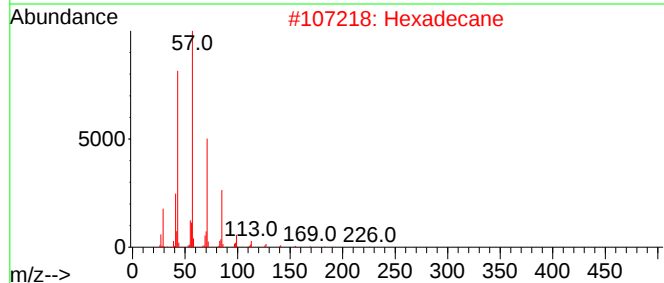
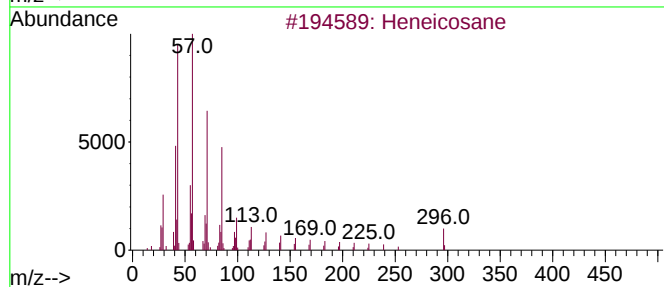
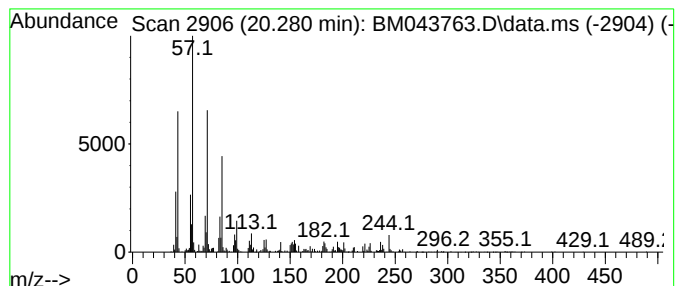
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.280	3.33 ng/ul	680193	Chrysene-d12		21.533	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	95
2	Hexadecane		226	C16H34	000544-76-3	95
3	Tricosane		324	C23H48	000638-67-5	95
4	Hexadecane		226	C16H34	000544-76-3	91
5	Hexadecane		226	C16H34	000544-76-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043763.D
Acq On : 05 Jan 2024 16:49
Operator : MA/JU
Sample : 05952-07ME
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF90ME

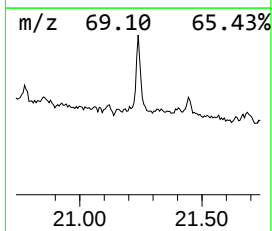
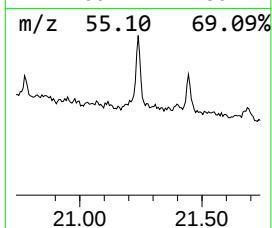
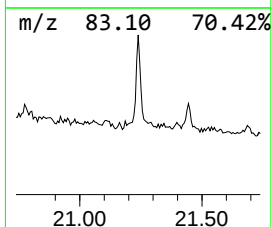
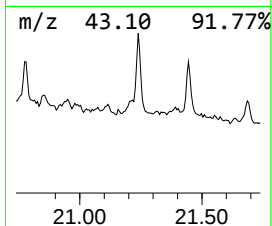
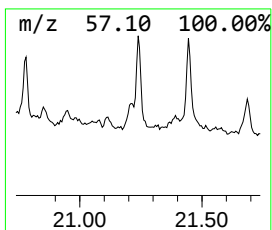
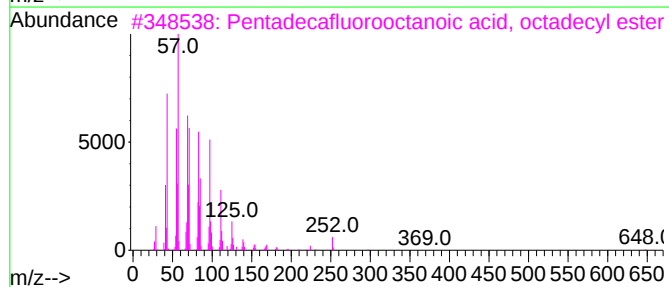
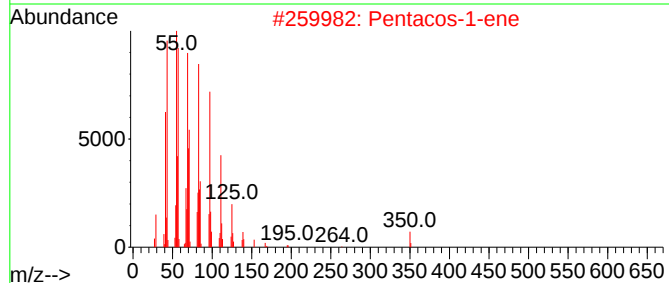
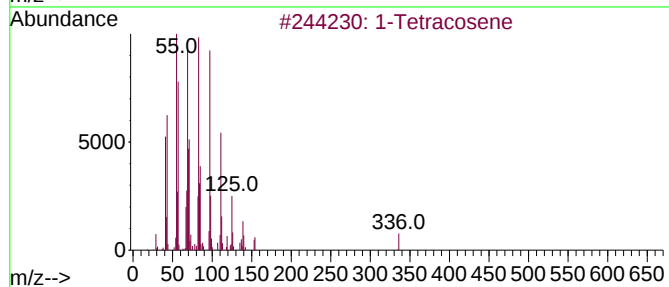
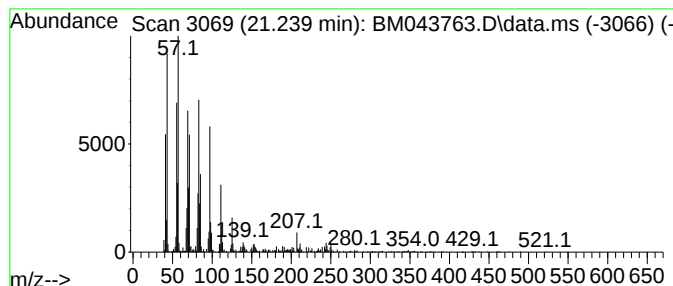
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.239	6.60 ng/ul	1350740	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tetracosene	336	C24H48	010192-32-2	97
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
4			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	90
5			Nonacos-1-ene	406	C29H58	018835-35-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043763.D
 Acq On : 05 Jan 2024 16:49
 Operator : MA/JU
 Sample : 05952-07ME
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF90ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 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
(DEL) Alkane: S...	14.486	3.1	ng/ul	462557	3	14.598	2934880	20.0
Naphthalene, 1...	14.992	3.1	ng/ul	460439	3	14.598	2934880	20.0
Naphthalene, 2...	15.069	2.0	ng/ul	300986	3	14.598	2934880	20.0
(DEL) Alkane: S...	15.433	5.7	ng/ul	830436	3	14.598	2934880	20.0
(DEL) Alkane: S...	15.845	8.8	ng/ul	1291550	3	14.598	2934880	20.0
Dibenzofuran, 4...	15.939	3.9	ng/ul	574044	3	14.598	2934880	20.0
(DEL) Alkane: S...	16.321	23.2	ng/ul	4474630	4	17.351	3853190	20.0
Anisole, 2,3,4...	16.704	6.3	ng/ul	1204830	4	17.351	3853190	20.0
unknown-01	16.880	2.5	ng/ul	489227	4	17.351	3853190	20.0
(DEL) Alkane: S...	17.098	14.9	ng/ul	2871720	4	17.351	3853190	20.0
(DEL) Alkane: S...	17.145	32.4	ng/ul	6242350	4	17.351	3853190	20.0
(DEL) Alkane: S...	17.756	7.3	ng/ul	1410860	4	17.351	3853190	20.0
(DEL) Alkane: S...	17.839	20.6	ng/ul	3967670	4	17.351	3853190	20.0
2-Methyldibenzo...	17.933	13.5	ng/ul	2595650	4	17.351	3853190	20.0
Phenanthrene, 1...	18.233	24.8	ng/ul	4770310	4	17.351	3853190	20.0
Phenanthrene, 2...	18.274	23.5	ng/ul	4533000	4	17.351	3853190	20.0
1H-Cyclopropa[1...	18.415	6.2	ng/ul	1186420	4	17.351	3853190	20.0
(DEL) Alkane: S...	18.539	21.8	ng/ul	4198870	4	17.351	3853190	20.0
Benzo[c]cinnoli...	18.621	12.0	ng/ul	2309940	4	17.351	3853190	20.0
Phenanthrene, 3...	18.974	7.9	ng/ul	1525540	4	17.351	3853190	20.0
Phenanthrene, 1...	19.021	5.4	ng/ul	1041790	4	17.351	3853190	20.0
di-p-Tolylacety...	19.062	16.1	ng/ul	3104910	4	17.351	3853190	20.0
Phenanthrene, 2...	19.145	13.7	ng/ul	2643920	4	17.351	3853190	20.0
3-Phenylquinox...	19.239	5.2	ng/ul	1010200	4	17.351	3853190	20.0
1-Nitro-9-methy...	19.462	6.9	ng/ul	1406120	5	21.533	4090760	20.0
(DEL) Alkane: S...	20.280	3.3	ng/ul	680193	5	21.533	4090760	20.0
(DEL) Alkane: S...	21.239	6.6	ng/ul	1350740	5	21.533	4090760	20.0