

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
 Data File : BM043335.D
 Acq On : 14 Dec 2023 12:02
 Operator : MA/JU
 Sample : 05690-06DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCLH5DL

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-BM120723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043335.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.834	106	110	123	rBV	65351	129789	2.03%	0.178%
2	7.157	666	675	687	rVB2	271646	560470	8.76%	0.769%
3	7.334	699	705	713	rVB	193171	319365	4.99%	0.438%
4	7.534	732	739	744	rVV3	378394	744291	11.64%	1.022%
5	7.593	744	749	758	rVB	210314	338331	5.29%	0.464%
6	7.781	774	781	791	rBV	717846	1198649	18.74%	1.645%
7	7.998	811	818	827	rBV	1508537	2447615	38.27%	3.360%
8	8.704	930	938	945	rBV	293769	479071	7.49%	0.658%
9	8.828	951	959	968	rBV	417469	692255	10.82%	0.950%
10	8.951	975	980	986	rVB	31453	53946	0.84%	0.074%
11	9.169	1010	1017	1025	rBV	239888	407946	6.38%	0.560%
12	9.886	1133	1139	1147	rBV	165540	275919	4.31%	0.379%
13	10.028	1155	1163	1172	rBV	59004	126563	1.98%	0.174%
14	10.251	1195	1201	1207	rVB	36889	60777	0.95%	0.083%
15	10.422	1223	1230	1239	rVB	306990	528924	8.27%	0.726%
16	10.804	1285	1295	1301	rBV	2192895	3809213	59.56%	5.229%
17	10.851	1301	1303	1311	rVV	79709	120764	1.89%	0.166%
18	10.951	1311	1320	1330	rVV	110137	228246	3.57%	0.313%
19	11.351	1381	1388	1394	rBV5	39887	77490	1.21%	0.106%
20	12.169	1519	1527	1539	rBV4	19025	54342	0.85%	0.075%
21	12.457	1569	1576	1583	rBV	84319	146202	2.29%	0.201%
22	12.675	1603	1613	1622	rVB	64330	120029	1.88%	0.165%
23	13.522	1752	1757	1766	rBV	38225	78451	1.23%	0.108%
24	13.792	1795	1803	1805	rBV5	38495	73806	1.15%	0.101%
25	13.945	1823	1829	1834	rBV	66342	122469	1.91%	0.168%
26	14.039	1840	1845	1851	rVB	574485	789239	12.34%	1.083%
27	14.186	1865	1870	1873	rVV2	32505	54246	0.85%	0.074%
28	14.316	1886	1892	1905	rVB	619502	1035706	16.19%	1.422%
29	14.621	1935	1944	1950	rVV	3332850	4887940	76.43%	6.710%
30	14.686	1950	1955	1962	rVV	481506	737476	11.53%	1.012%
31	14.821	1972	1978	1982	rBV2	175257	315792	4.94%	0.433%
32	14.857	1982	1984	1990	rVV3	72007	96778	1.51%	0.133%
33	14.933	1990	1997	2001	rVV4	31314	62417	0.98%	0.086%
34	15.021	2007	2012	2019	rVV	223337	340461	5.32%	0.467%
35	15.404	2069	2077	2081	rBV8	22382	52094	0.81%	0.072%
36	15.610	2106	2112	2117	rVV	831144	1229765	19.23%	1.688%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
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37	15.668	2117	2122	2127	rVV	412038	621499	9.72%	0.853%
38	15.727	2128	2132	2136	rVV3	80098	136379	2.13%	0.187%
39	15.792	2139	2143	2146	rVV2	65267	106594	1.67%	0.146%
40	15.827	2146	2149	2153	rVV2	58078	93205	1.46%	0.128%

41	15.874	2153	2157	2161	rVV4	37950	70228	1.10%	0.096%
42	15.957	2167	2171	2179	rVB	76122	127778	2.00%	0.175%
43	16.104	2188	2196	2204	rVB	67782	156124	2.44%	0.214%
44	16.333	2227	2235	2244	rBV6	62761	168569	2.64%	0.231%
45	16.668	2281	2292	2297	rBV3	73298	169621	2.65%	0.233%

46	16.721	2297	2301	2305	rVV2	35220	56330	0.88%	0.077%
47	16.821	2314	2318	2322	rVB3	35719	56490	0.88%	0.078%
48	16.898	2324	2331	2333	rVV3	29892	55975	0.88%	0.077%
49	17.015	2348	2351	2358	rVB5	40258	66846	1.05%	0.092%
50	17.098	2359	2365	2366	rVV2	41528	66016	1.03%	0.091%

51	17.121	2367	2369	2374	rVV2	51496	82601	1.29%	0.113%
52	17.180	2374	2379	2388	rVB	144134	235665	3.68%	0.324%
53	17.362	2404	2410	2414	rBV	3716966	5318027	83.15%	7.300%
54	17.404	2414	2417	2422	rVV	1951560	2632774	41.17%	3.614%
55	17.462	2422	2427	2430	rVV2	814581	1255962	19.64%	1.724%

56	17.498	2430	2433	2438	rVV	472296	631830	9.88%	0.867%
57	17.551	2439	2442	2447	rVB2	37149	66590	1.04%	0.091%
58	17.768	2474	2479	2491	rBV	96977	202929	3.17%	0.279%
59	18.239	2554	2559	2564	rBV2	173101	393913	6.16%	0.541%
60	18.286	2564	2567	2575	rVB	257507	353276	5.52%	0.485%

61	18.368	2577	2581	2585	rVV	97690	142193	2.22%	0.195%
62	18.433	2586	2592	2596	rVV2	600588	975439	15.25%	1.339%
63	18.474	2596	2599	2605	rVB2	359529	514957	8.05%	0.707%
64	18.645	2623	2628	2638	rBV4	321374	488901	7.64%	0.671%
65	18.739	2640	2644	2648	rVV	153332	206505	3.23%	0.283%

66	18.780	2648	2651	2655	rVV3	49750	74875	1.17%	0.103%
67	18.833	2656	2660	2669	rVB3	169096	274871	4.30%	0.377%
68	18.980	2680	2685	2690	rBV	566037	700043	10.95%	0.961%
69	19.062	2696	2699	2703	rVB3	75792	94822	1.48%	0.130%
70	19.109	2703	2707	2710	rBV3	133529	187328	2.93%	0.257%

71	19.151	2710	2714	2718	rVB	90264	129882	2.03%	0.178%
72	19.309	2734	2741	2745	rBV6	85810	214193	3.35%	0.294%
73	19.415	2753	2759	2764	rBV	3411998	4515442	70.60%	6.198%
74	19.474	2764	2769	2773	rVB	2351524	3021471	47.24%	4.148%
75	19.686	2802	2805	2809	rVB2	168118	203352	3.18%	0.279%

76	19.745	2810	2815	2817	rVV3	1438246	2030605	31.75%	2.787%
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77	19.774	2817	2820	2829	rVV	2886726	3833368	59.94%	5.262%
78	19.845	2829	2832	2839	rVB2	121585	181865	2.84%	0.250%
79	19.956	2847	2851	2855	rVB	122028	168277	2.63%	0.231%
80	20.098	2871	2875	2882	rVB5	147607	310596	4.86%	0.426%
81	20.233	2895	2898	2899	rBV3	77797	89641	1.40%	0.123%
82	20.268	2900	2904	2908	rVB2	428316	562123	8.79%	0.772%
83	20.368	2917	2921	2924	rBV	420806	501882	7.85%	0.689%
84	20.421	2925	2930	2934	rVB3	200216	358938	5.61%	0.493%
85	20.562	2951	2954	2957	rBV	123318	146008	2.28%	0.200%
86	20.609	2959	2962	2966	rVB3	121928	175085	2.74%	0.240%
87	20.733	2980	2983	2987	rVB2	126737	140522	2.20%	0.193%
88	21.180	3056	3059	3062	rBV	185627	219641	3.43%	0.302%
89	21.215	3062	3065	3068	rVV	193592	241035	3.77%	0.331%
90	21.256	3068	3072	3077	rVB3	307662	482266	7.54%	0.662%
91	21.533	3112	3119	3123	rBV2	3731204	6395578	100.00%	8.779%
92	21.568	3123	3125	3133	rVB	1094329	1345188	21.03%	1.847%
93	21.692	3143	3146	3154	rVB2	184870	326221	5.10%	0.448%
94	22.174	3225	3228	3235	rVB4	195923	311890	4.88%	0.428%
95	23.215	3399	3405	3409	rBV	771911	1774779	27.75%	2.436%
96	23.733	3489	3493	3498	rBV	371533	638274	9.98%	0.876%
97	23.791	3499	3503	3507	rVV2	383847	624934	9.77%	0.858%
98	23.839	3508	3511	3516	rVB	376395	611337	9.56%	0.839%
99	23.950	3524	3530	3537	rVB	1775284	3371846	52.72%	4.629%
100	24.638	3644	3647	3655	rVB	194975	339204	5.30%	0.466%

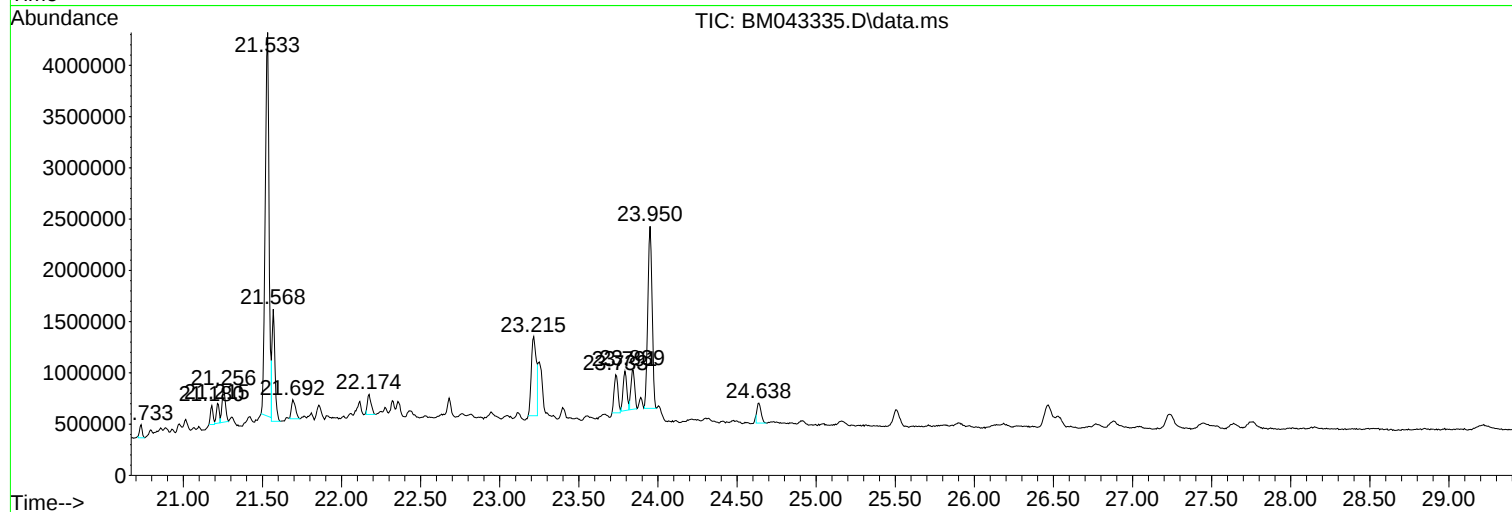
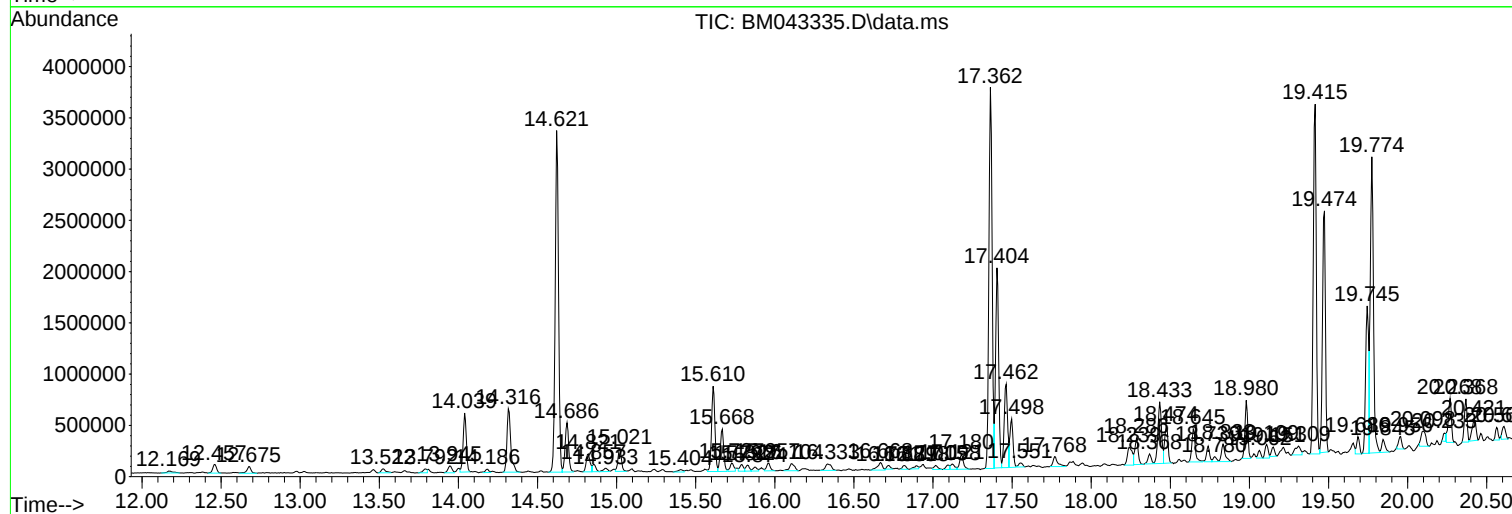
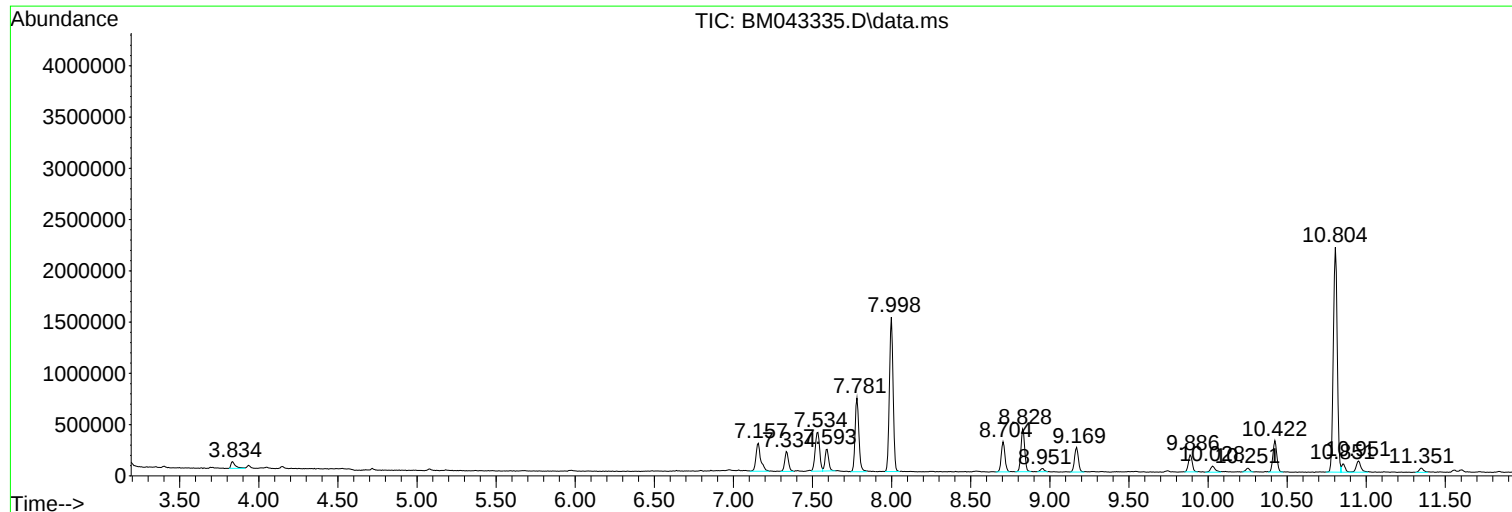
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ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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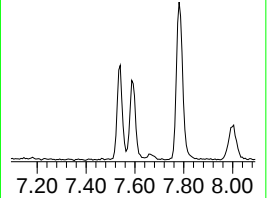
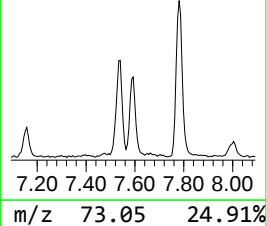
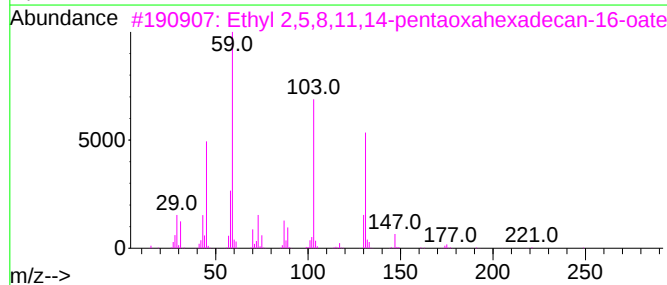
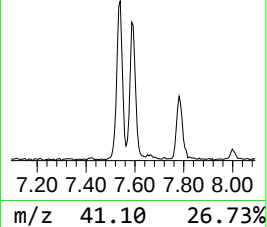
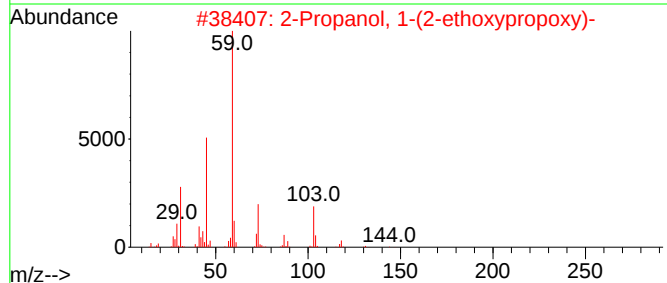
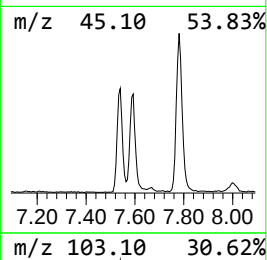
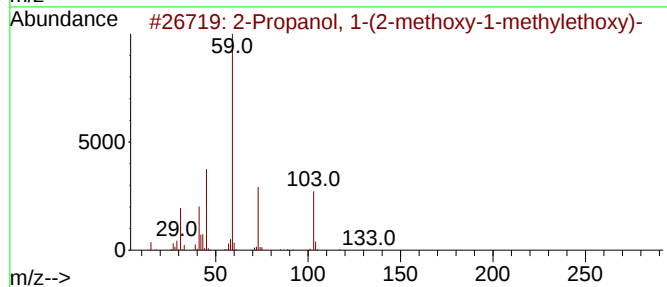
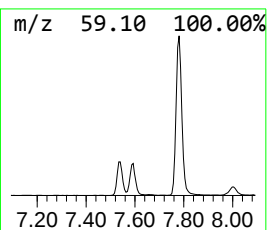
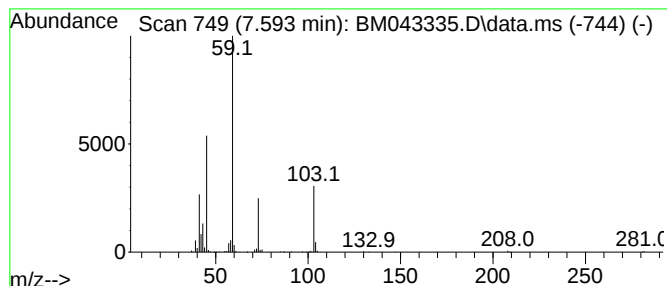
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.593	2.76 ng/ul	338331	1,4-Dichlorobenzene-d4	7.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72		
2	2-Propanol, 1-(2-ethoxypropoxy)-	162	C8H18O3	010143-32-5	64		
3	Ethyl 2,5,8,11,14-pentaoxahexade...	294	C13H26O7	1000366-80-7	56		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	53		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	53		



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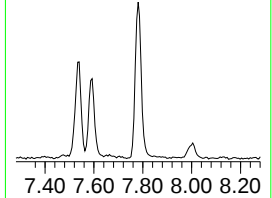
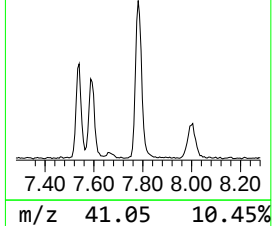
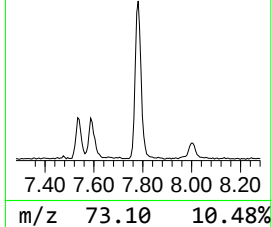
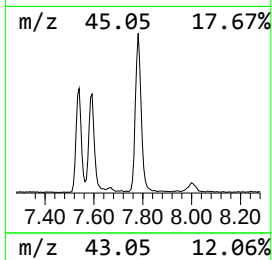
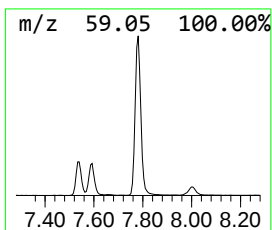
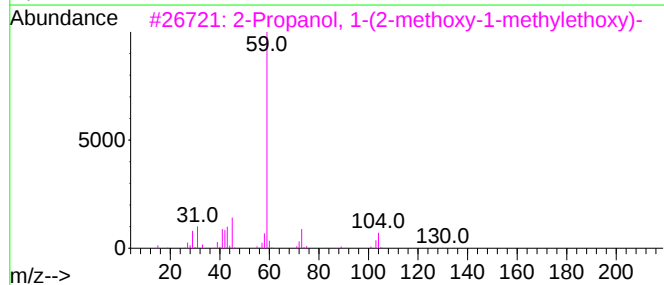
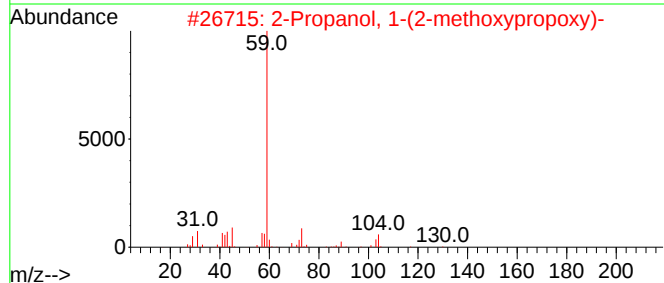
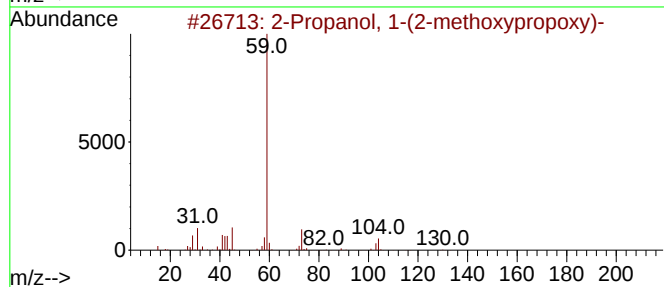
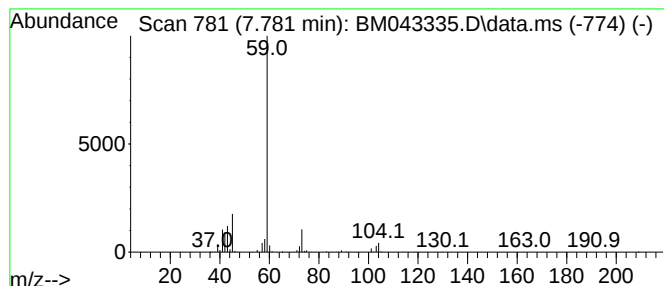
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propanol, 1-(2-methoxypro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.781	9.79 ng/ul	1198650	1,4-Dichlorobenzene-d4	7.998

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	83	
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	74	
3	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64	
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64	
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	45	



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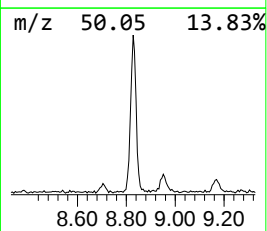
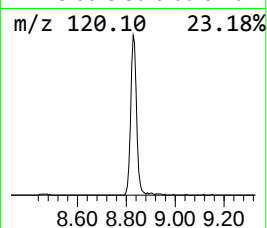
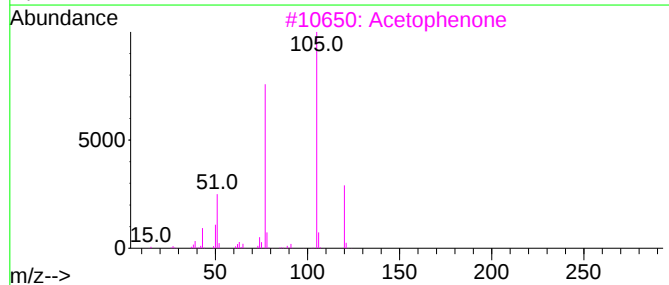
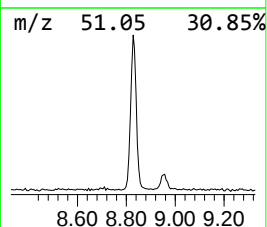
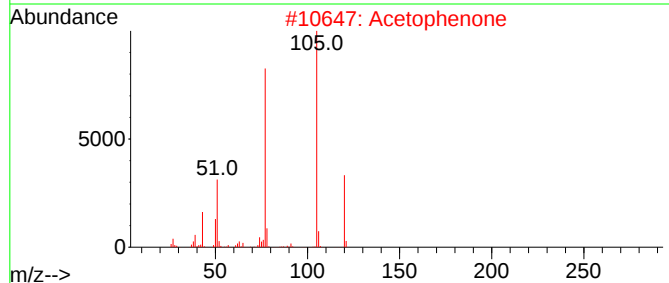
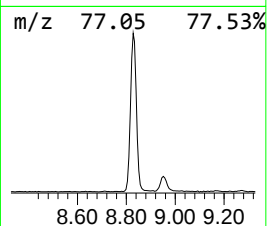
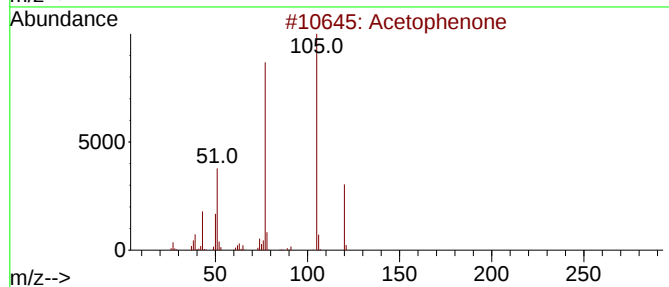
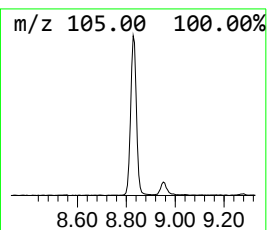
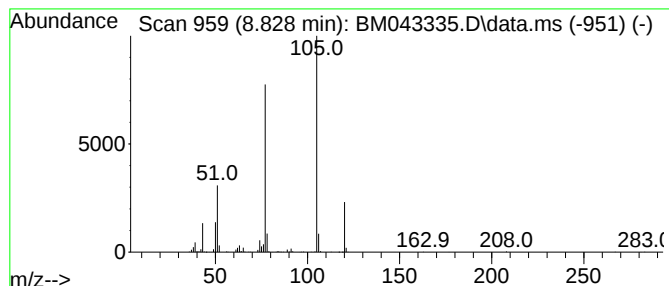
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Aceto phenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.828	5.66 ng/ul	692255	1,4-Dichlorobenzene-d4	7.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	97
2	Acetophenone			120	C8H8O	000098-86-2	95
3	Acetophenone			120	C8H8O	000098-86-2	94
4	Acetophenone			120	C8H8O	000098-86-2	94
5	Acetophenone			120	C8H8O	000098-86-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

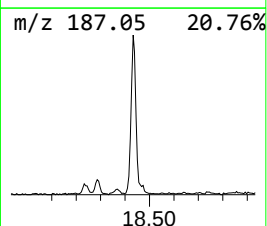
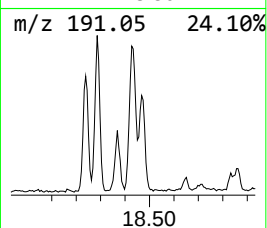
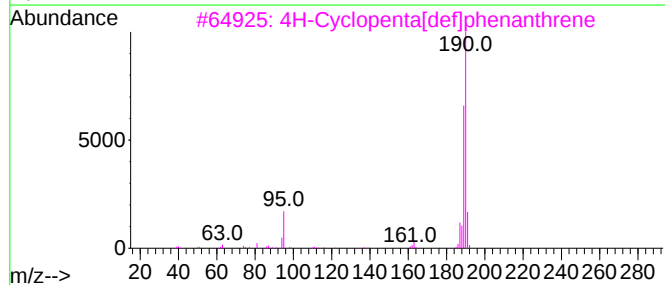
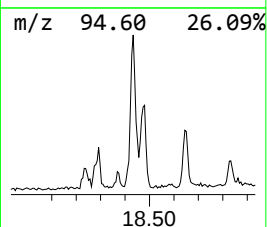
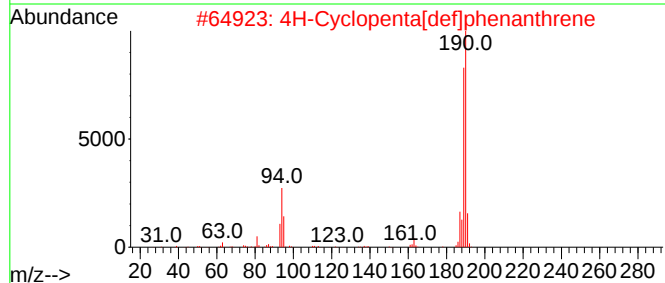
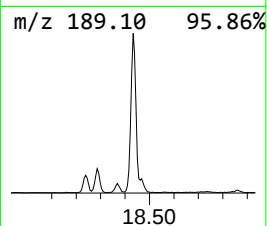
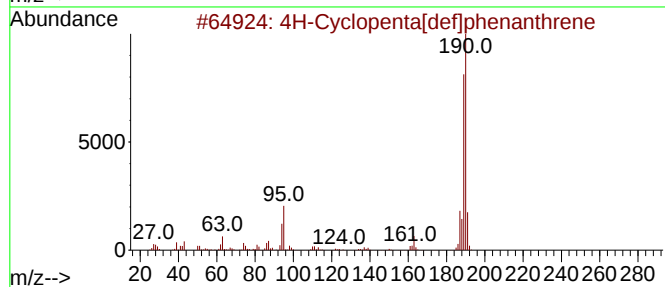
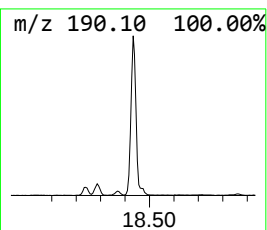
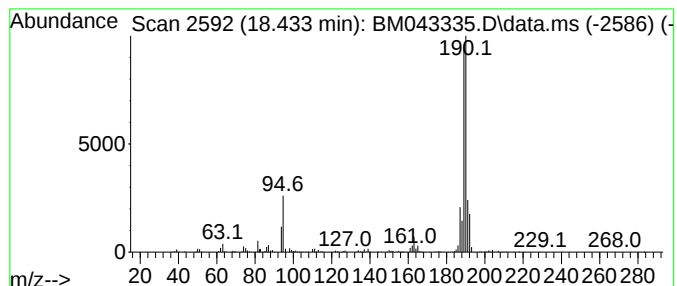
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.433	3.67 ng/ul	975439	Phenanthrene-d10	17.363	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	80
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
5	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLH5DL

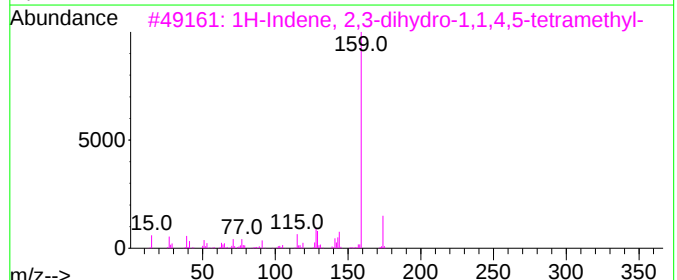
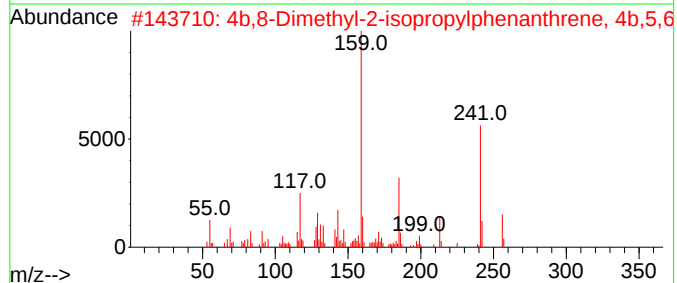
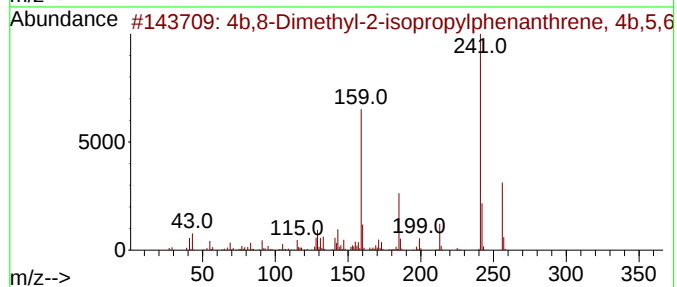
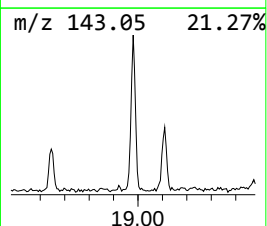
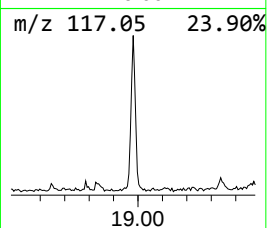
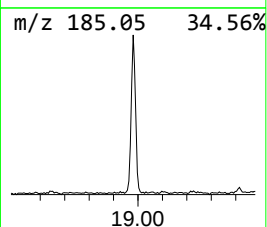
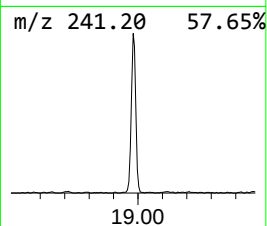
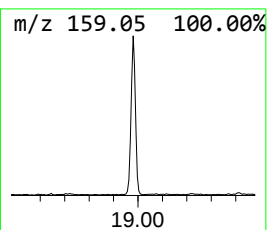
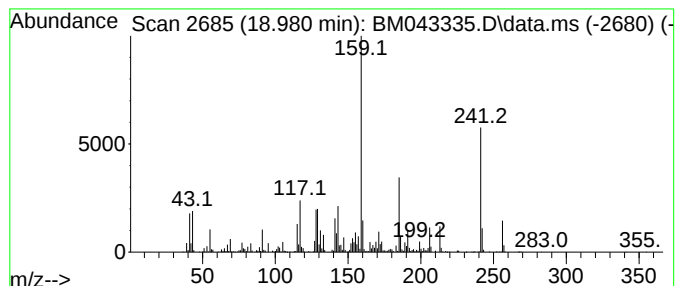
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4b,8-Dimethyl-2-isopropylph... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.980	2.63 ng/ul	700043	Phenanthrene-d10	17.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	98
2			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	93
3			1H-Indene, 2,3-dihydro-1,1,4,5-t...	174	C13H18	016204-57-2	38
4			2(1H)-Quinolinone, hydrazone	159	C9H9N3	015793-77-8	38
5			2-Phenylpropionic acid, 4'-(2-me...	218	C14H18O2	1010454-94-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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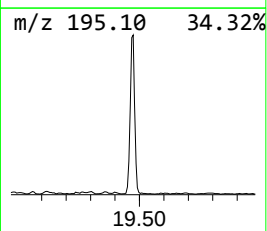
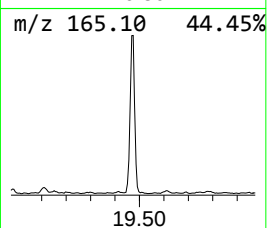
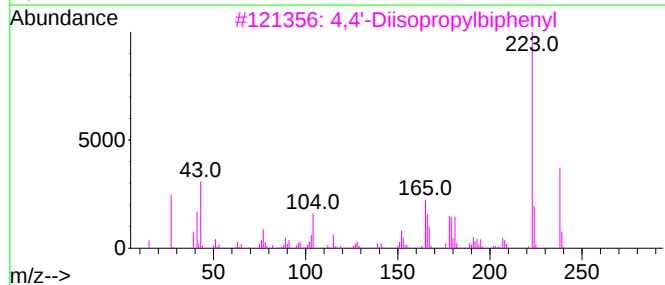
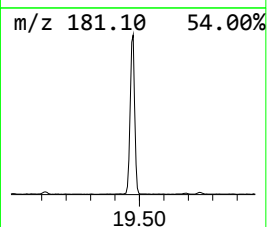
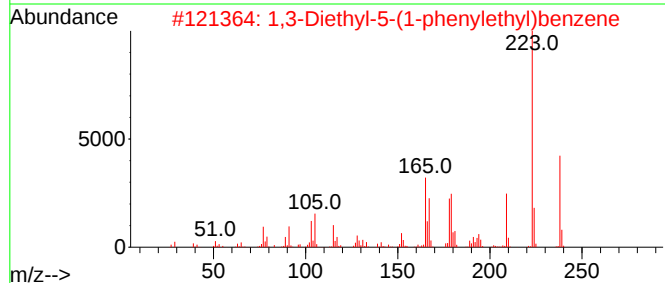
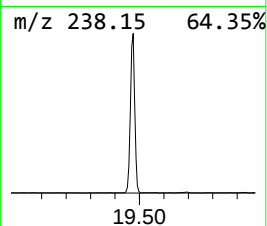
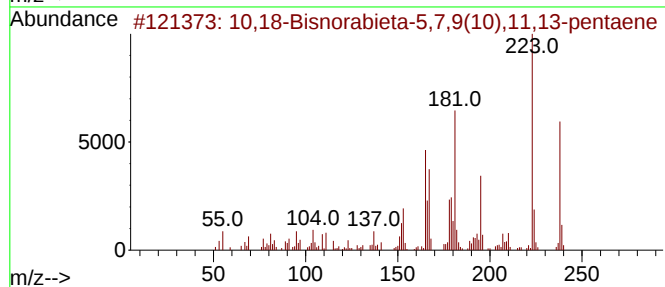
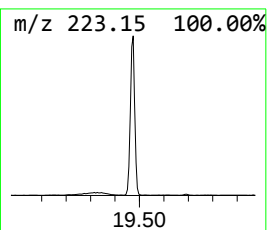
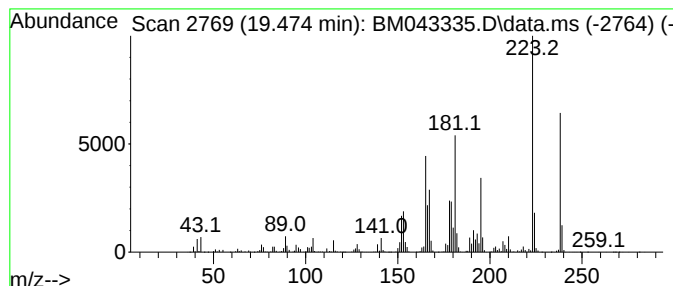
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.474	9.45 ng/ul	3021470	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	72		
3	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	55		
4	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	50		
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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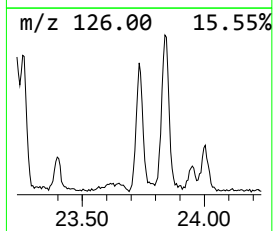
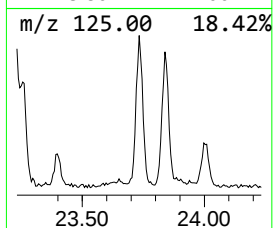
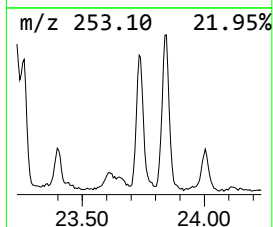
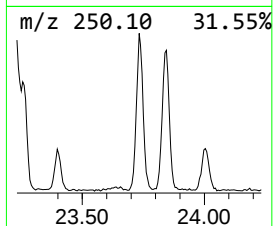
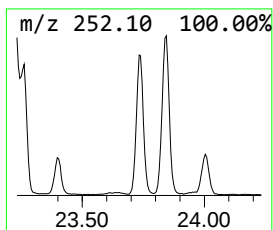
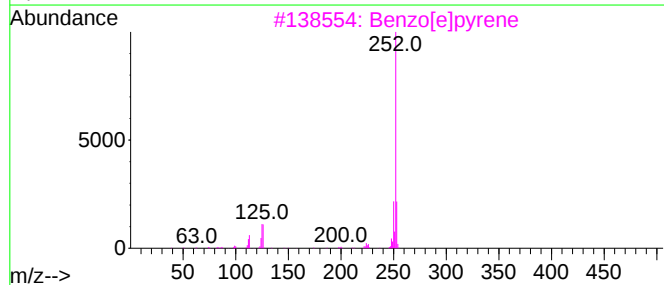
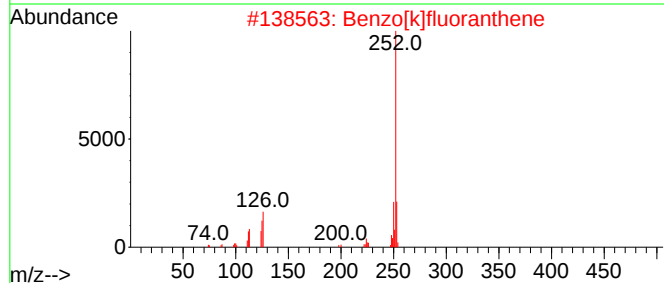
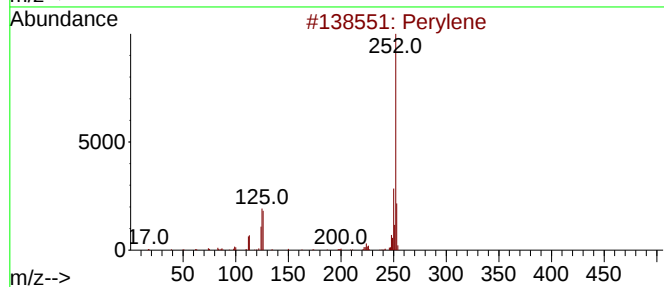
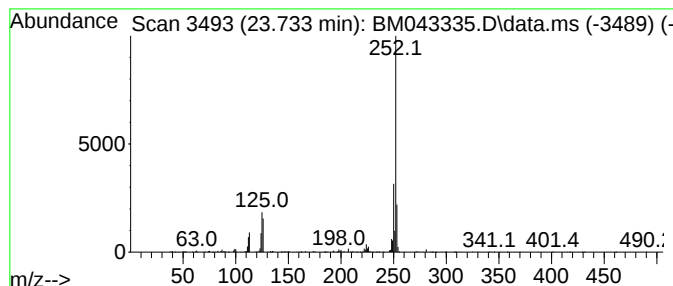
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Perylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.733	3.79 ng/ul	638274	Perylene-d12	23.950

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	99
3			Benzo[e]pyrene	252	C20H12	000192-97-2	98
4			Benzo[e]pyrene	252	C20H12	000192-97-2	98
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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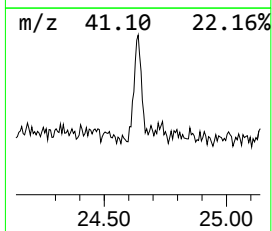
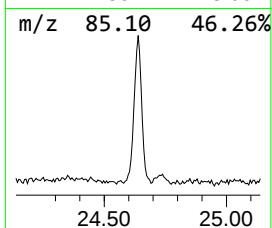
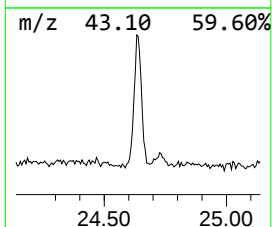
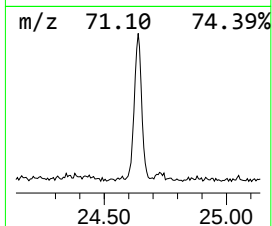
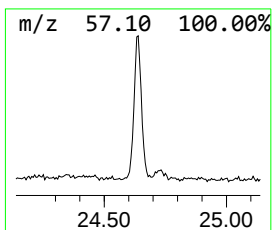
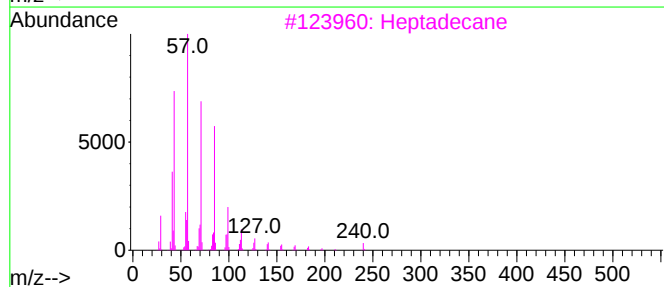
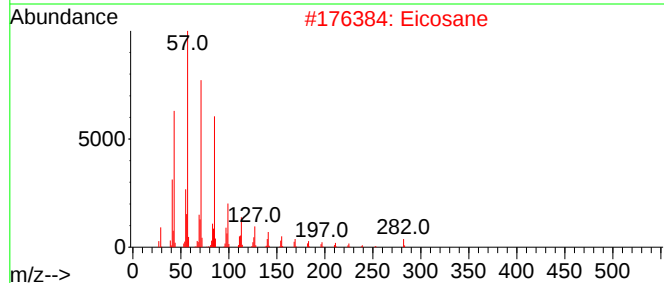
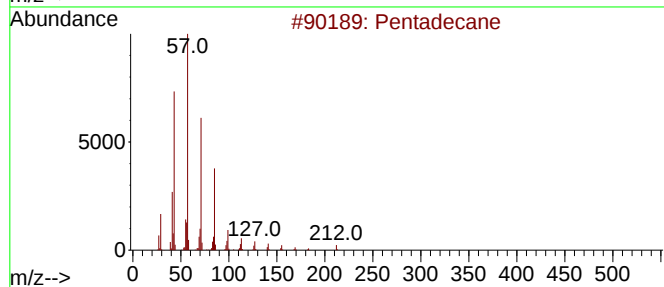
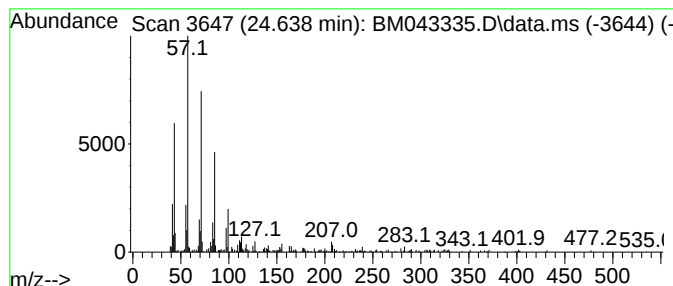
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.639	2.01 ng/ul	339204	Perylene-d12	23.950

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	90
2		Eicosane	282	C20H42	000112-95-8	87
3		Heptadecane	240	C17H36	000629-78-7	87
4		Pentacosane	352	C25H52	000629-99-2	86
5		Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043335.D
Acq On : 14 Dec 2023 12:02
Operator : MA/JU
Sample : 05690-06DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLH5DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Propanol, 1-(...	7.781	9.8	ng/ul	1198650	1	7.998	2447620	20.0
Aceto phenone	8.828	5.7	ng/ul	692255	1	7.998	2447620	20.0
4H-Cyclopenta[d...	18.433	3.7	ng/ul	975439	4	17.363	5318030	20.0
4b,8-Dimethyl-2...	18.980	2.6	ng/ul	700043	4	17.363	5318030	20.0
10,18-Bisnorabi...	19.474	9.4	ng/ul	3021470	5	21.533	6395580	20.0
Perylene	23.733	3.8	ng/ul	638274	6	23.950	3371850	20.0
(DEL) Alkane: S...	24.639	2.0	ng/ul	339204	6	23.950	3371850	20.0