

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
 Data File : BM047744.D
 Acq On : 01 Oct 2024 15:43
 Operator : RC/JU
 Sample : P4018-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDCA2

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-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047744.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.834	478	484	493	rVB2	3220835	4992188	5.08%	1.065%
2	6.134	531	535	542	rVB	1147305	1807096	1.84%	0.385%
3	6.316	559	566	572	rBV4	976853	1864362	1.90%	0.398%
4	6.381	572	577	581	rBV	1030745	1442847	1.47%	0.308%
5	6.457	581	590	593	rVV3	825645	1764970	1.80%	0.376%
6	6.716	626	634	640	rVB4	580384	1305182	1.33%	0.278%
7	6.816	646	651	655	rVB2	1420999	2381075	2.42%	0.508%
8	6.869	655	660	663	rBV	1787870	2498239	2.54%	0.533%
9	6.904	663	666	671	rVB	1429514	2093693	2.13%	0.446%
10	6.975	671	678	684	rBV2	3044568	5833353	5.93%	1.244%
11	7.040	684	689	694	rVB	1462778	2128695	2.17%	0.454%
12	7.140	701	706	711	rBV	821941	1305349	1.33%	0.278%
13	7.204	711	717	721	rVV2	1146219	1921727	1.95%	0.410%
14	7.334	731	739	745	rVV2	2975337	5497569	5.59%	1.172%
15	7.445	752	758	766	rVB2	9374561	16660223	16.95%	3.553%
16	7.739	800	808	813	rVV4	657674	2013419	2.05%	0.429%
17	7.804	813	819	825	rVV2	2531604	4492900	4.57%	0.958%
18	7.887	825	833	838	rVV	8459543	14190507	14.44%	3.026%
19	7.934	838	841	846	rVV2	995284	1687940	1.72%	0.360%
20	8.010	846	854	855	rVV2	867791	1681675	1.71%	0.359%
21	8.039	855	859	868	rVV3	2137364	5353290	5.45%	1.142%
22	8.110	868	871	878	rVV3	829253	1404638	1.43%	0.300%
23	8.245	889	894	899	rBV2	668891	1343284	1.37%	0.286%
24	8.351	907	912	918	rVV2	1744725	3576283	3.64%	0.763%
25	8.410	918	922	925	rVV	1258084	1692869	1.72%	0.361%
26	8.457	925	930	931	rVV2	1584112	2214068	2.25%	0.472%
27	8.481	931	934	937	rVV	2038155	3137324	3.19%	0.669%
28	8.510	937	939	946	rVB2	1479582	2107491	2.14%	0.449%
29	8.586	946	952	955	rBV	1757659	2708112	2.75%	0.578%
30	8.616	955	957	965	rVB	958876	1489989	1.52%	0.318%
31	8.769	978	983	987	rBV	840579	1257472	1.28%	0.268%
32	8.816	987	991	999	rVB	936524	1508708	1.53%	0.322%
33	8.922	1000	1009	1012	rBV2	1191185	2482806	2.53%	0.529%
34	8.963	1012	1016	1020	rVV	832549	1365701	1.39%	0.291%
35	9.051	1021	1031	1036	rVB	7431346	11393449	11.59%	2.430%
36	9.175	1040	1052	1058	rVB7	1733085	6128766	6.23%	1.307%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

37	9.245	1058	1064	1069	rBV2	1392582	2539203	2.58%	0.541%
38	9.686	1132	1139	1145	rBV4	1049437	2532919	2.58%	0.540%
39	9.757	1145	1151	1156	rBV3	1144190	2120741	2.16%	0.452%
40	10.228	1227	1231	1239	rVB2	1214780	2258294	2.30%	0.482%
41	10.492	1268	1276	1283	rBV	2093120	4089110	4.16%	0.872%
42	10.592	1286	1293	1300	rVV3	4533889	8980329	9.14%	1.915%
43	10.733	1311	1317	1322	rBV3	2566456	4885048	4.97%	1.042%
44	10.904	1339	1346	1354	rVB	838868	1540597	1.57%	0.329%
45	10.986	1354	1360	1368	rBV	1833221	3242092	3.30%	0.691%
46	11.298	1405	1413	1419	rBV3	933199	1942631	1.98%	0.414%
47	11.639	1466	1471	1474	rBV	1056583	1705307	1.73%	0.364%
48	11.704	1474	1482	1490	rVB	3380311	6395081	6.51%	1.364%
49	11.875	1503	1511	1520	rVB	1688742	2780179	2.83%	0.593%
50	12.033	1532	1538	1542	rBV	2417815	4137400	4.21%	0.882%
51	12.069	1542	1544	1549	rVB	1638602	2120614	2.16%	0.452%
52	12.133	1549	1555	1559	rBV	914305	1655240	1.68%	0.353%
53	12.280	1573	1580	1589	rVB2	2744739	5468922	5.56%	1.166%
54	12.680	1642	1648	1660	rBV2	1777962	3953865	4.02%	0.843%
55	13.280	1745	1750	1757	rVB	3017156	4357119	4.43%	0.929%
56	13.716	1819	1824	1828	rBV	909730	1350619	1.37%	0.288%
57	13.769	1828	1833	1836	rVV	976226	1687831	1.72%	0.360%
58	13.816	1836	1841	1844	rVV4	1331705	2190524	2.23%	0.467%
59	13.851	1844	1847	1854	rVB	1952300	2825971	2.87%	0.603%
60	14.080	1882	1886	1890	rVB	1701442	1978240	2.01%	0.422%
61	14.139	1891	1896	1900	rBV	1786305	2524406	2.57%	0.538%
62	14.363	1928	1934	1938	rBV	17303297	25924816	26.37%	5.529%
63	14.445	1938	1948	1953	rBV	2384354	3905048	3.97%	0.833%
64	14.592	1968	1973	1977	rBV	4699591	6020412	6.12%	1.284%
65	14.645	1977	1982	1993	rVB6	1199154	3490419	3.55%	0.744%
66	14.763	1994	2002	2005	rBV3	1554191	3409307	3.47%	0.727%
67	14.915	2022	2028	2032	rBV3	817521	1574442	1.60%	0.336%
68	14.974	2033	2038	2041	rVV2	745783	1405175	1.43%	0.300%
69	15.010	2041	2044	2047	rVV	902799	1328332	1.35%	0.283%
70	15.074	2047	2055	2060	rVV	10967881	17616270	17.92%	3.757%
71	15.133	2062	2065	2070	rVV	1616322	2247038	2.29%	0.479%
72	15.210	2073	2078	2085	rVV	5053779	7871846	8.01%	1.679%
73	15.304	2086	2094	2098	rVB2	3678551	6605255	6.72%	1.409%
74	15.439	2112	2117	2122	rVB	2686457	3839083	3.91%	0.819%
75	15.545	2129	2135	2142	rBV	2312924	3500428	3.56%	0.746%
76	15.710	2157	2163	2168	rBV	1142377	1803466	1.83%	0.385%

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77	15.798	2174	2178	2184	rVB2	1370344	2091792	2.13%	0.446%
78	16.074	2220	2225	2230	rVB	1999548	2716146	2.76%	0.579%
79	16.162	2235	2240	2243	rBV	3765131	5299932	5.39%	1.130%
80	16.268	2253	2258	2263	rBV	2780329	4038971	4.11%	0.861%
81	16.504	2294	2298	2302	rBV3	981242	1416908	1.44%	0.302%
82	16.874	2348	2361	2366	rVV2	33701315	98305986	100.00%	20.964%
83	16.951	2370	2374	2379	rVV	3128561	4186025	4.26%	0.893%
84	17.004	2379	2383	2393	rVB2	1492251	2634276	2.68%	0.562%
85	17.192	2409	2415	2419	rBV	3059332	4526474	4.60%	0.965%
86	17.286	2426	2431	2435	rVV	3019994	4178525	4.25%	0.891%
87	17.327	2435	2438	2443	rVB3	1020196	1264478	1.29%	0.270%
88	17.480	2459	2464	2471	rVB4	1055094	1966012	2.00%	0.419%
89	17.686	2494	2499	2504	rVB	2605766	3277295	3.33%	0.699%
90	17.856	2524	2528	2532	rVB	2203166	2555279	2.60%	0.545%
91	18.080	2562	2566	2573	rVB2	867979	1340567	1.36%	0.286%
92	18.374	2612	2616	2626	rBV	1946111	2577953	2.62%	0.550%
93	19.027	2719	2727	2731	rVB2	1944601	4005111	4.07%	0.854%
94	19.568	2813	2819	2826	rBV2	3504295	5987190	6.09%	1.277%
95	20.115	2908	2912	2927	rVB2	1114166	1782645	1.81%	0.380%
96	21.092	3073	3078	3092	rVB	1808799	2718214	2.77%	0.580%
97	21.403	3126	3131	3137	rBV2	2705260	4009661	4.08%	0.855%
98	22.174	3257	3262	3273	rVB3	709633	1288782	1.31%	0.275%
99	24.203	3599	3607	3619	rVB	1831669	4527348	4.61%	0.965%
100	24.409	3634	3642	3660	rVB	1965357	5701877	5.80%	1.216%

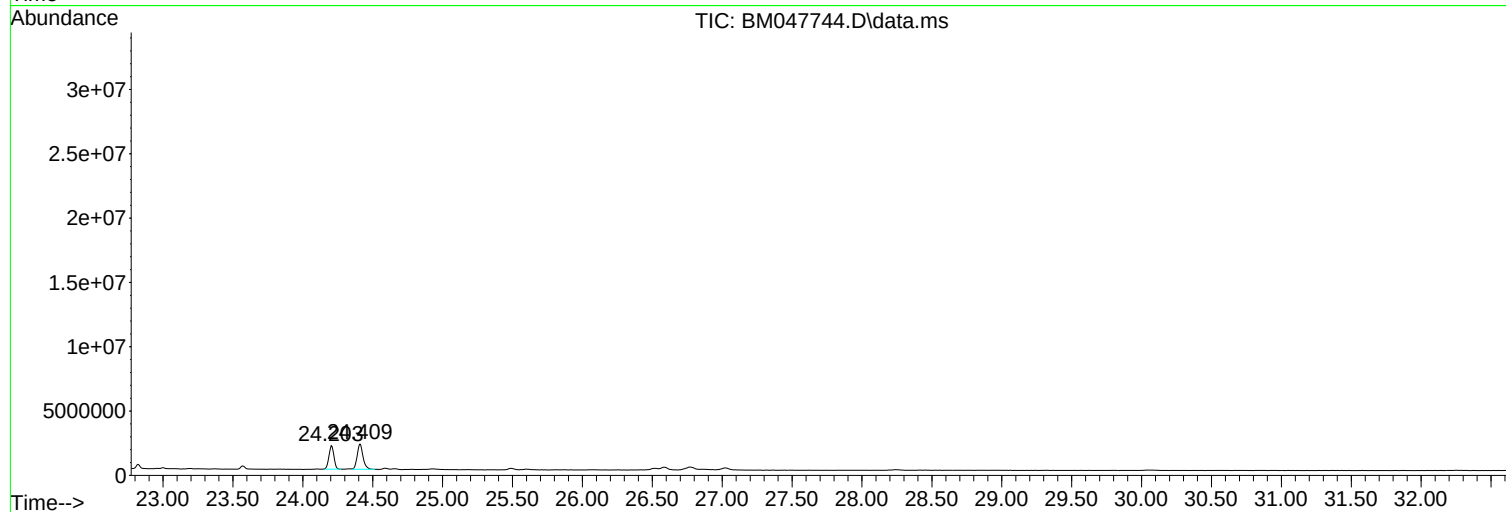
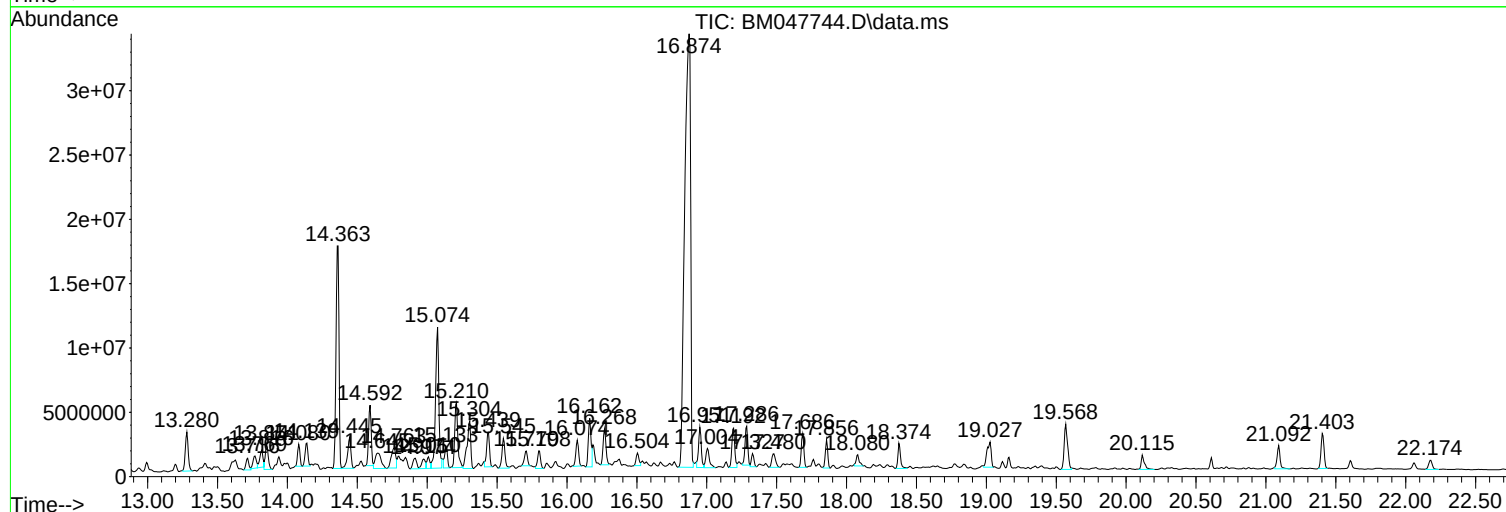
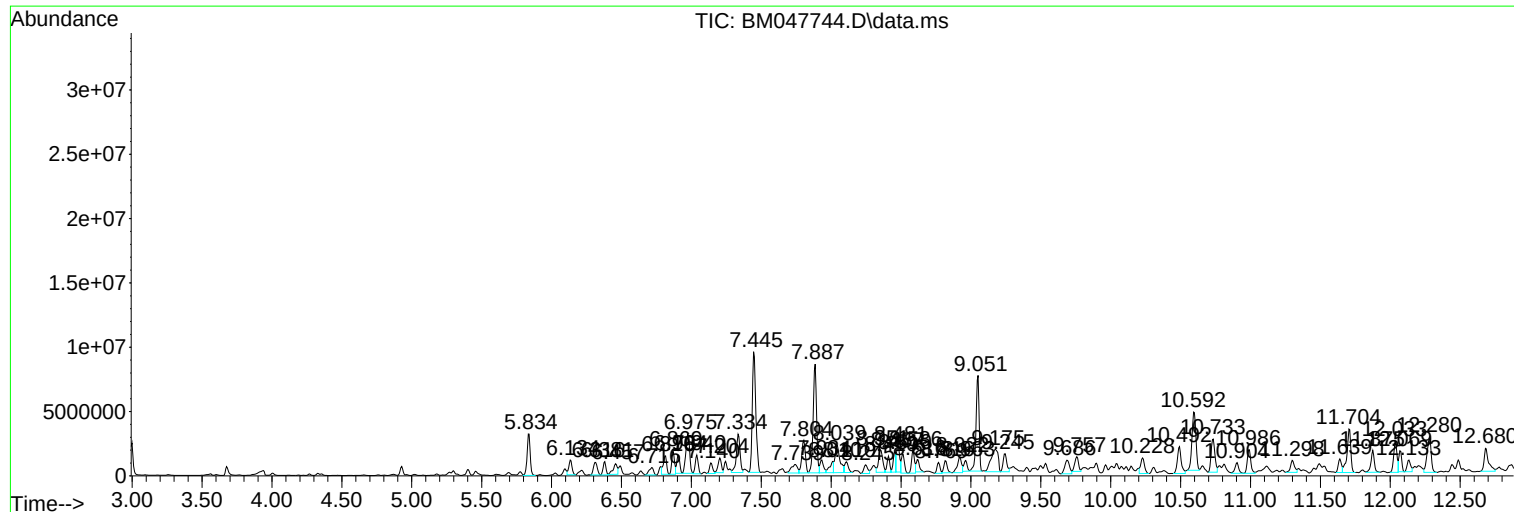
Sum of corrected areas: 468928325

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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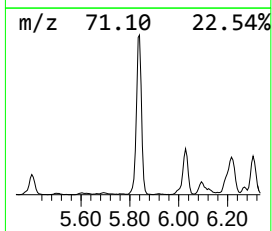
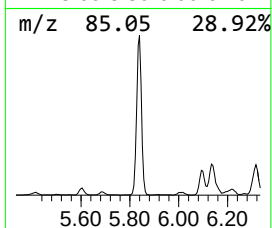
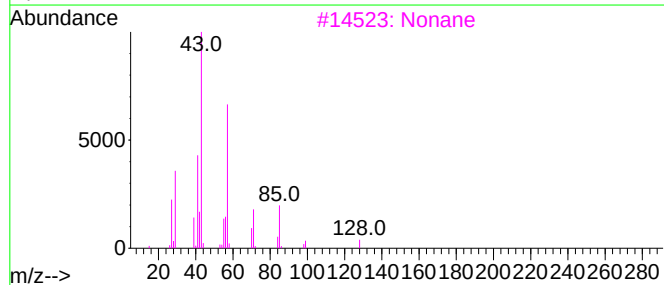
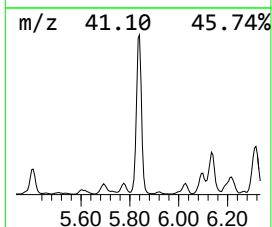
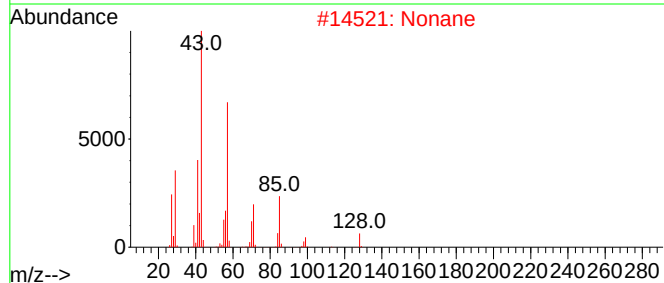
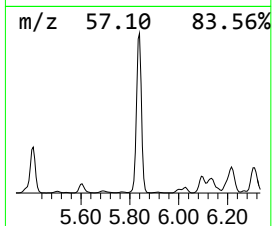
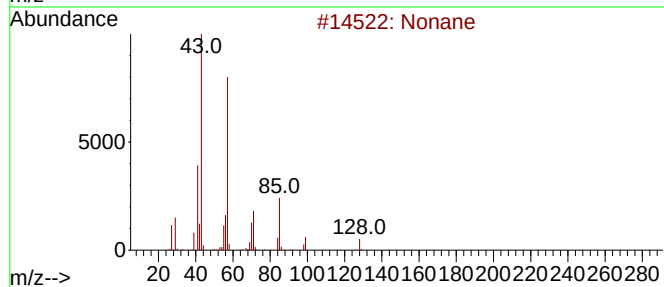
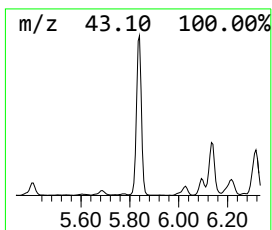
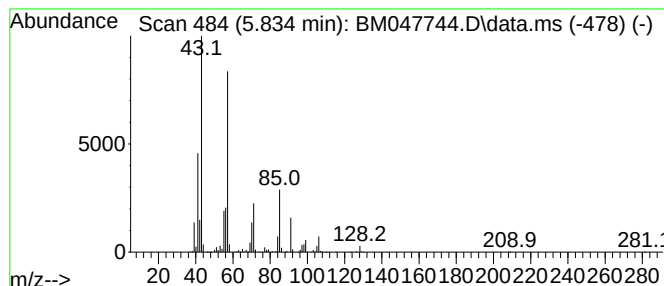
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.834	22.22 ng/ul	4992190	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	87
2			Nonane	128	C9H20	000111-84-2	80
3			Nonane	128	C9H20	000111-84-2	72
4			Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	53
5			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	53



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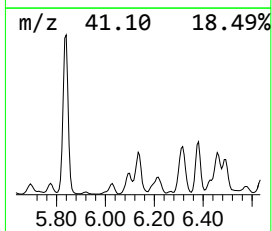
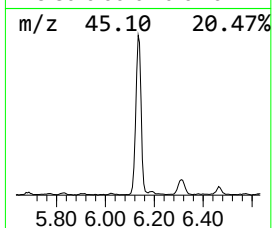
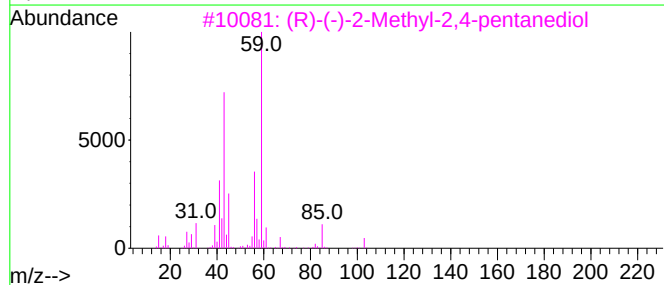
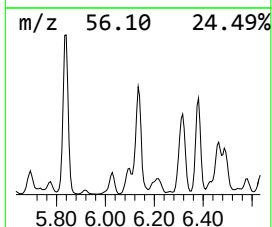
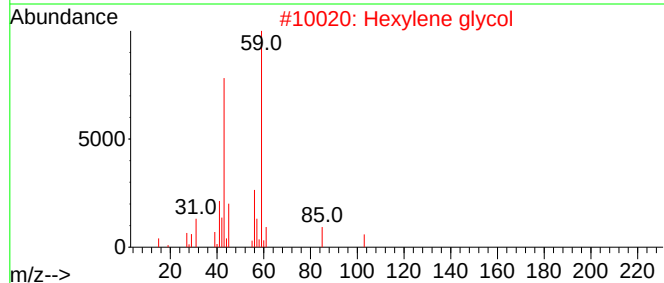
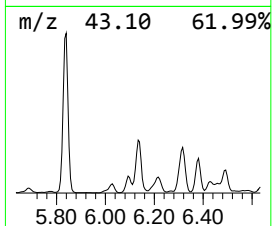
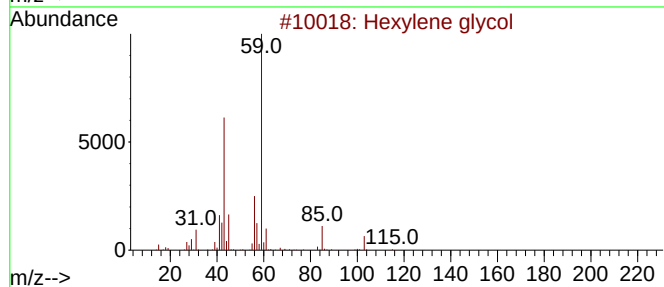
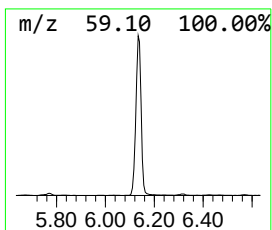
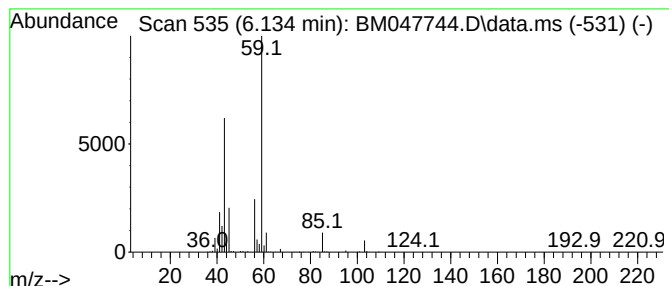
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.134	8.04 ng/ul	1807100	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	83
3			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	64
4			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
5			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59



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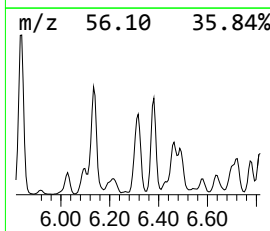
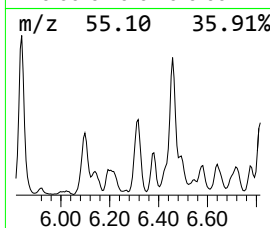
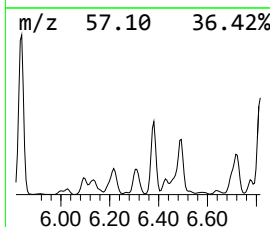
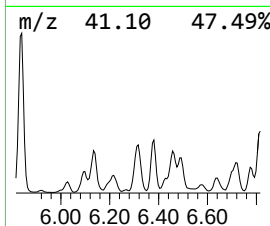
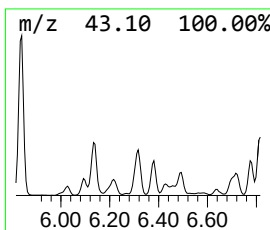
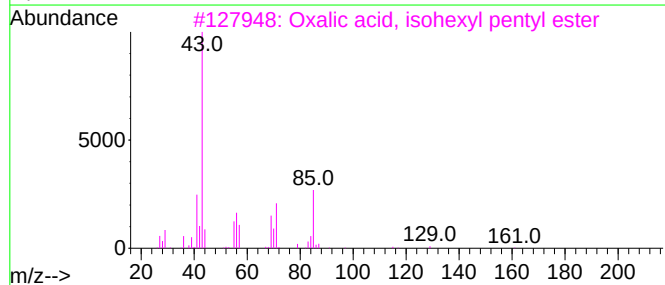
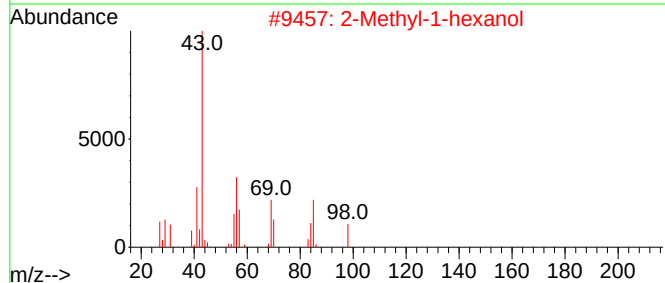
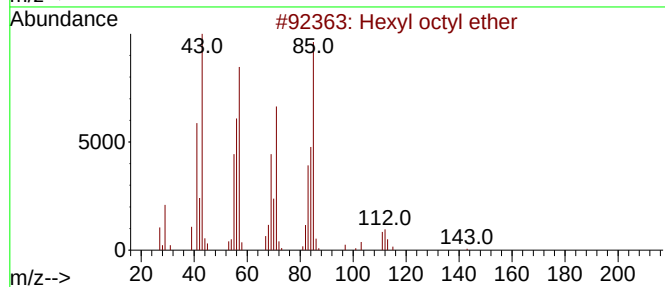
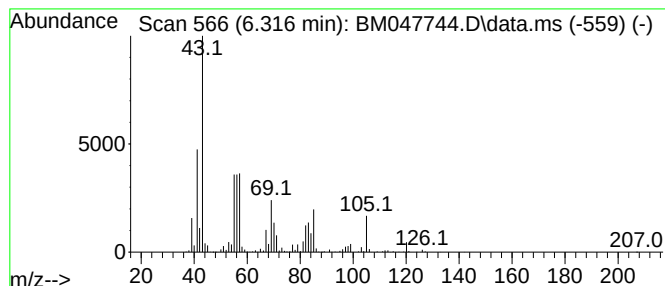
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexyl octyl ether Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.316	8.30 ng/ul	1864360	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexyl octyl ether	214	C14H30O	017071-54-4	50
2			2-Methyl-1-hexanol	116	C7H16O	1000427-67-0	43
3			Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	43
4			1-Pentanol, 2,3-dimethyl-	116	C7H16O	010143-23-4	38
5			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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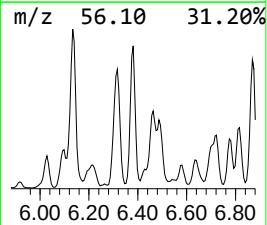
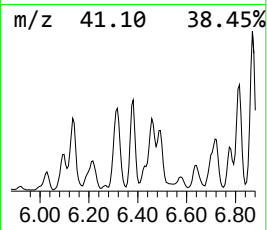
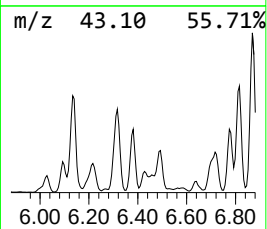
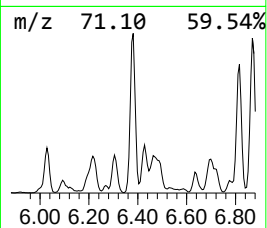
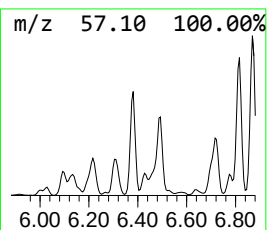
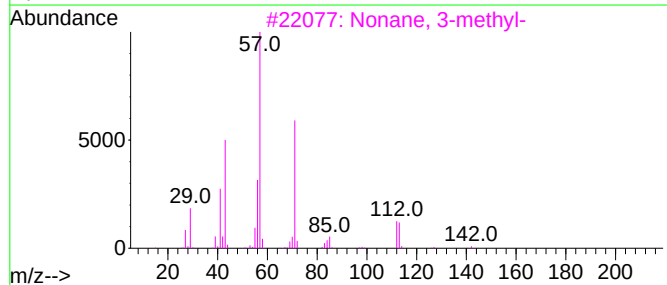
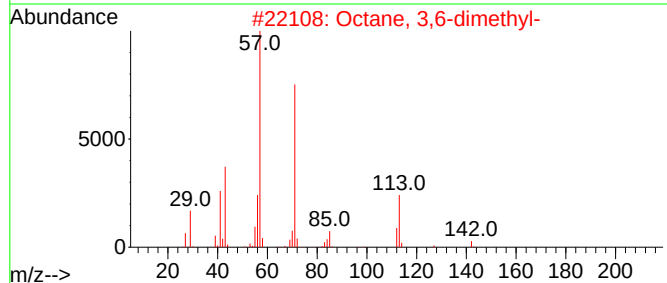
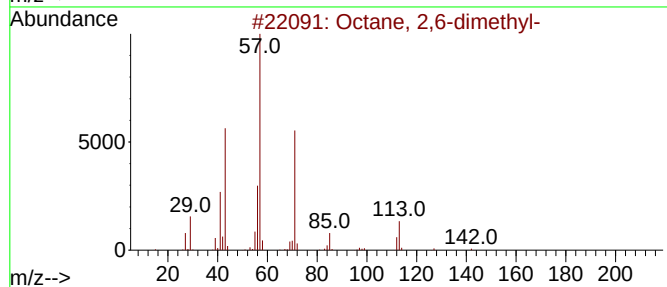
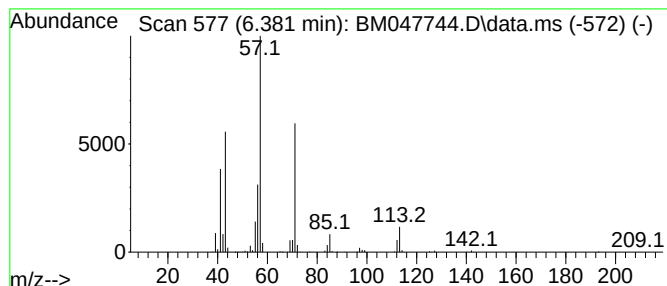
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.381	6.42 ng/ul	1442850	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	94
2	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	91
3	Nonane, 3-methyl-			142	C10H22	005911-04-6	90
4	Nonane, 3-methyl-			142	C10H22	005911-04-6	87
5	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
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Sample : P4018-04
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Instrument :
BNA_M
ClientSampleId :
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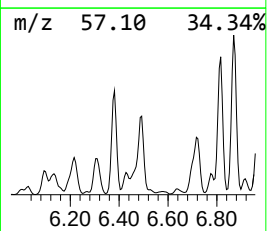
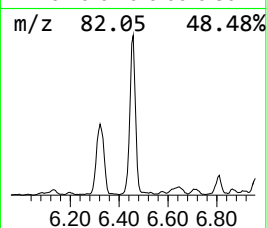
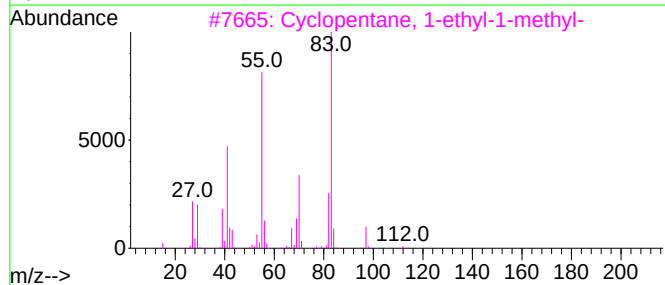
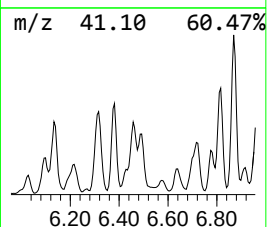
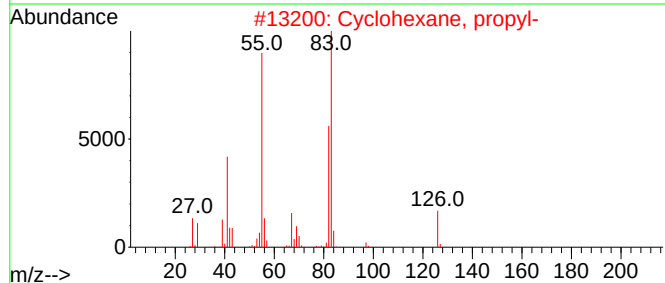
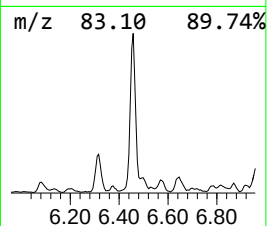
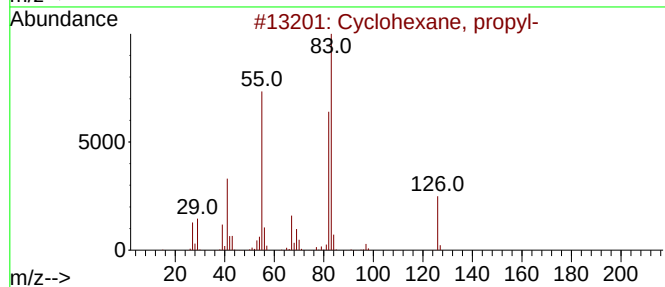
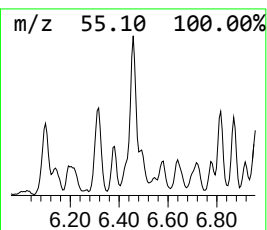
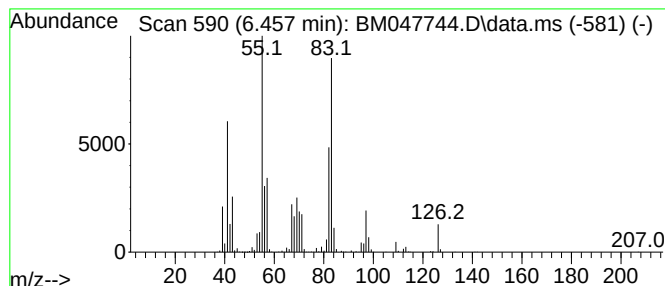
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic6.457 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.457	7.86 ng/ul	1764970	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	58
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
5			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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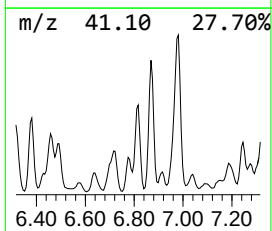
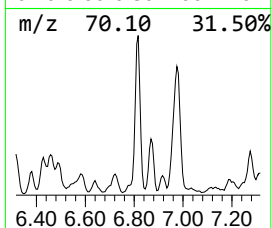
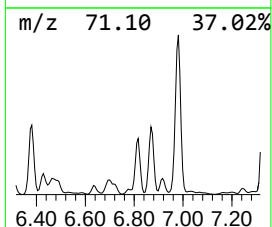
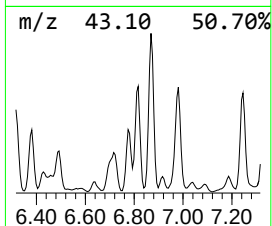
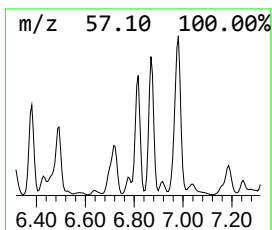
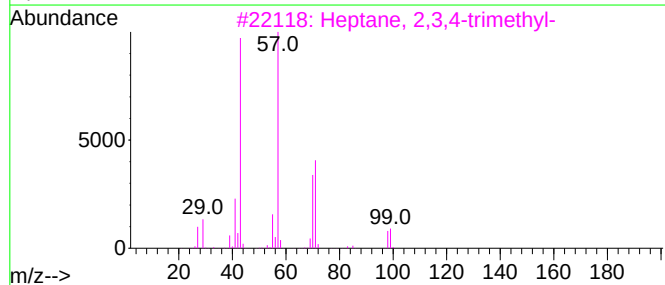
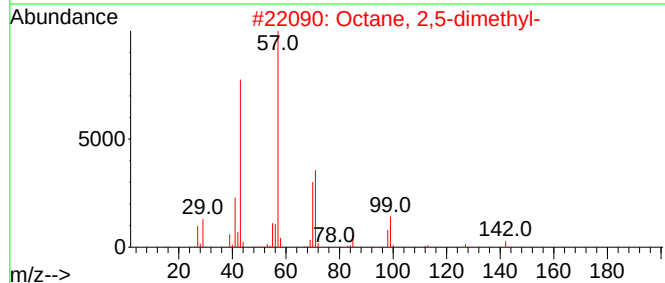
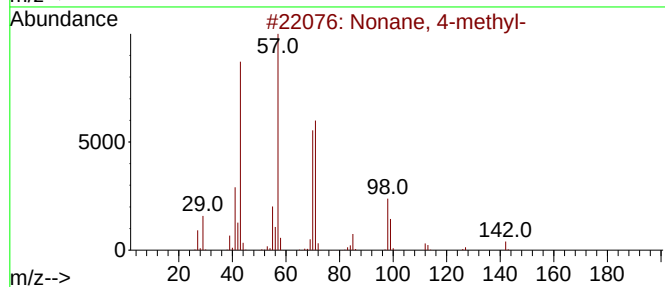
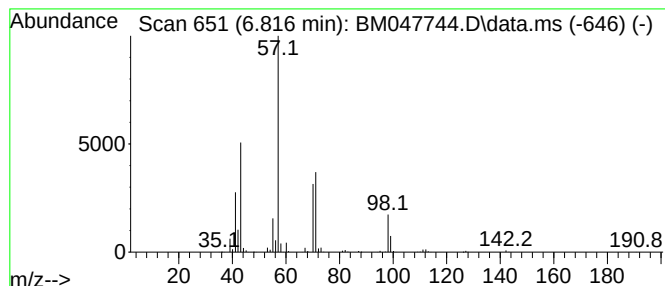
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.816	10.60 ng/ul	2381080	1,4-Dichlorobenzene-d4	7.810

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 4-methyl-	142	C10H22	017301-94-9	64
2		Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	59
3		Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	59
4		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
5		Nonane, 4-methyl-	142	C10H22	017301-94-9	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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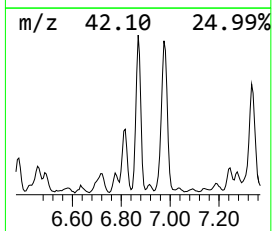
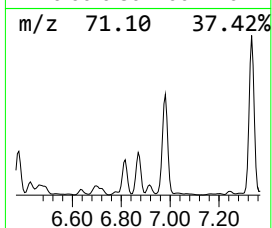
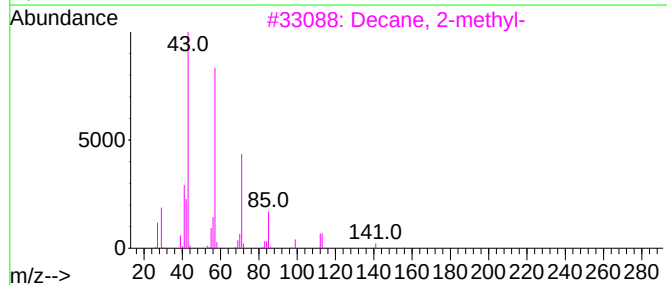
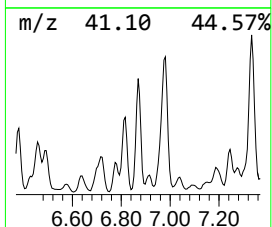
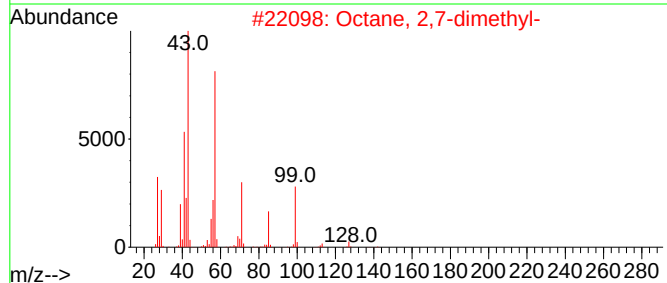
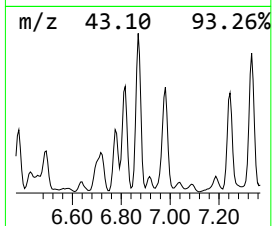
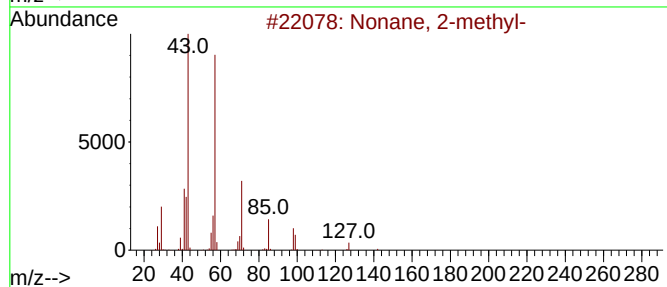
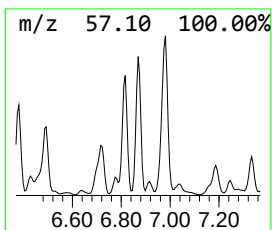
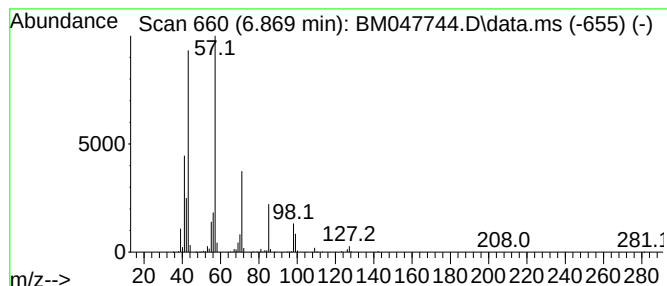
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.869	11.12 ng/ul	2498240	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	91
2			Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	72
3			Decane, 2-methyl-	156	C11H24	006975-98-0	64
4			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	59
5			Octane, 2-methyl-	128	C9H20	003221-61-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
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Instrument :
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ClientSampleId :
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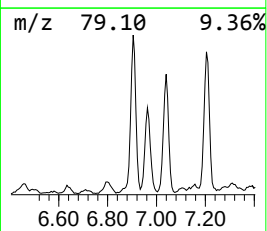
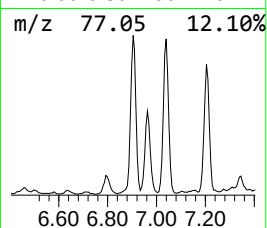
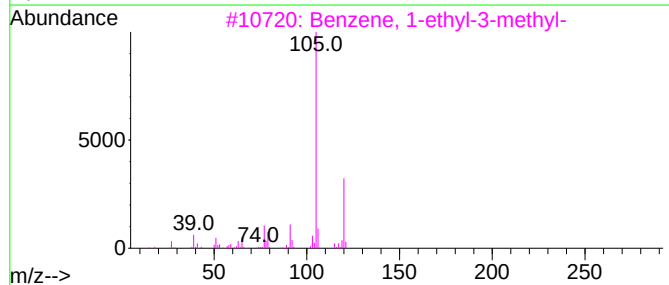
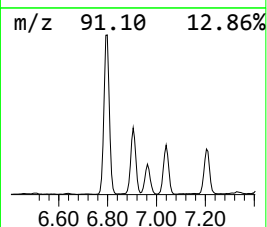
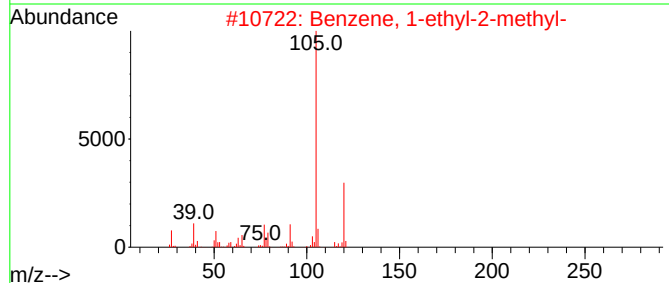
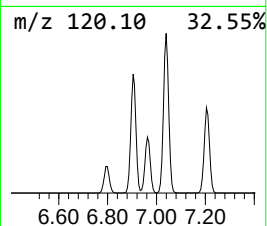
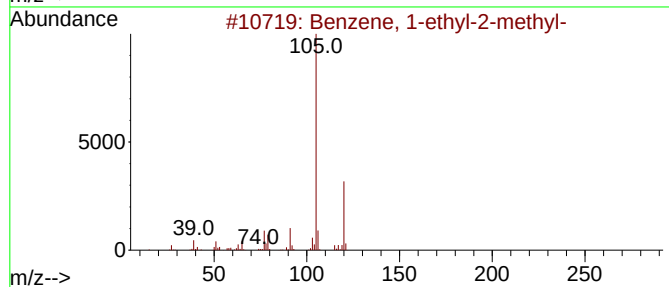
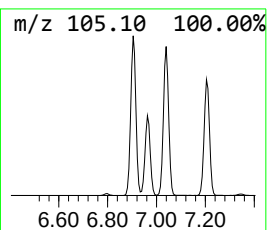
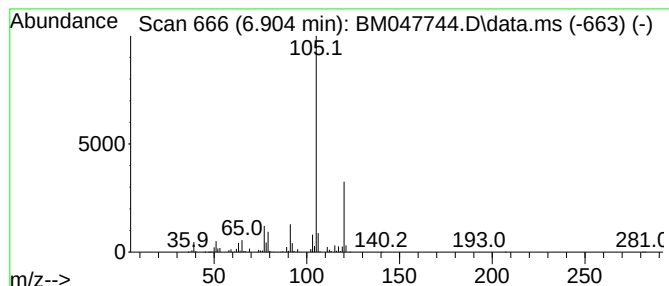
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.904	9.32 ng/ul	2093690	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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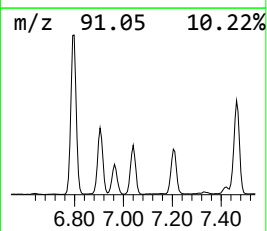
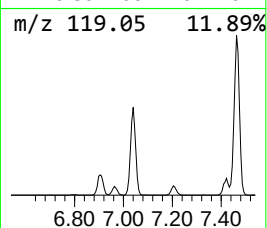
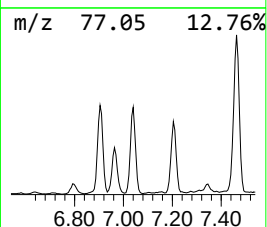
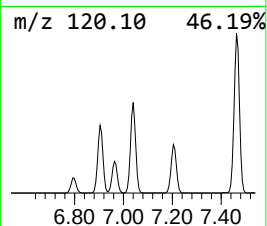
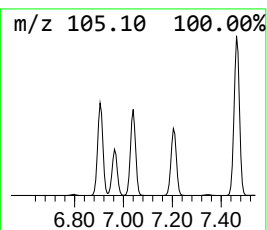
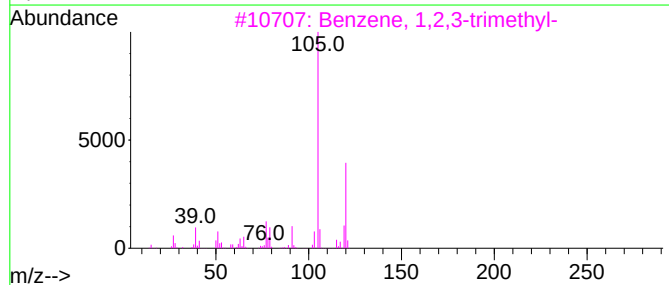
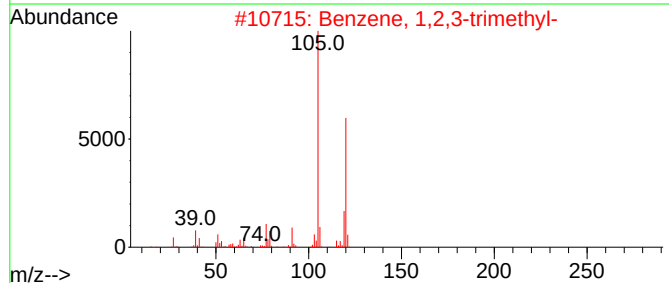
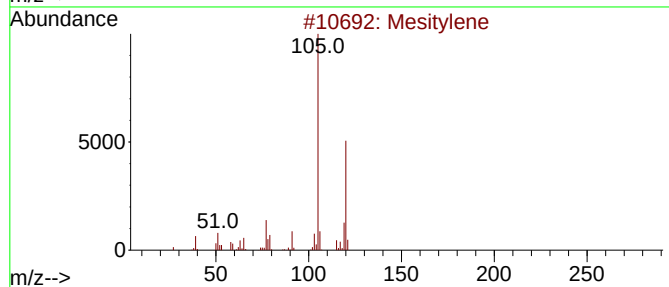
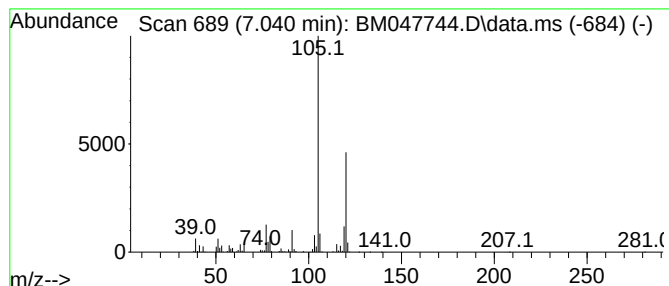
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Mesitylene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.040	9.48 ng/ul	2128700	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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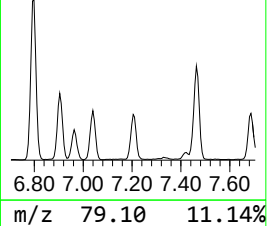
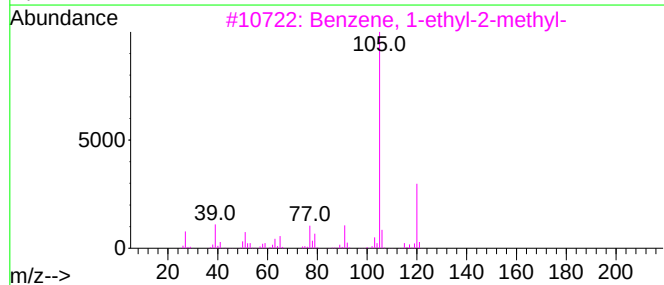
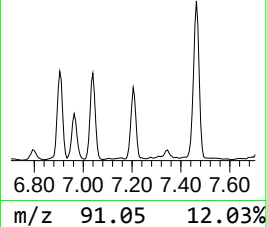
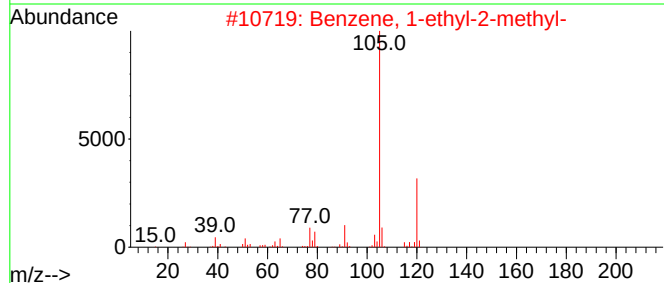
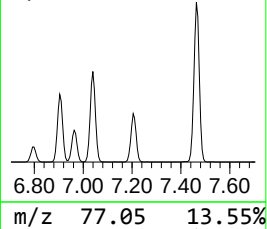
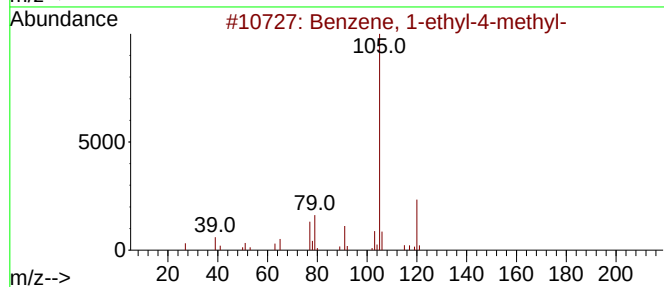
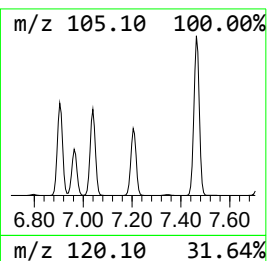
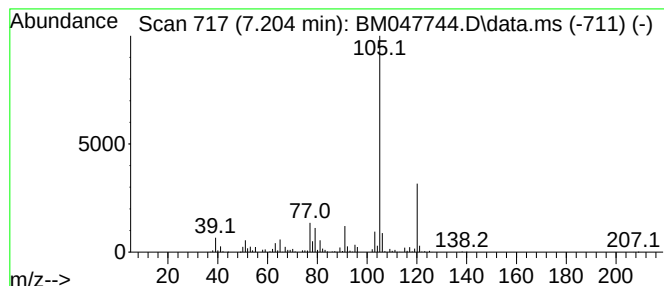
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1-ethyl-4-methyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.204	8.55 ng/ul	1921730	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Mesitylene	120	C9H12	000108-67-8	90
5			Mesitylene	120	C9H12	000108-67-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

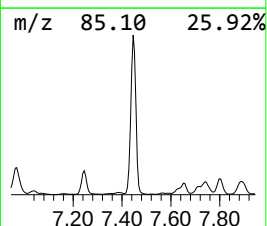
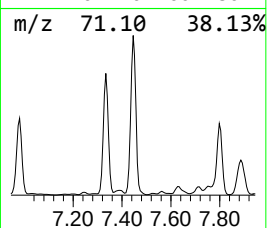
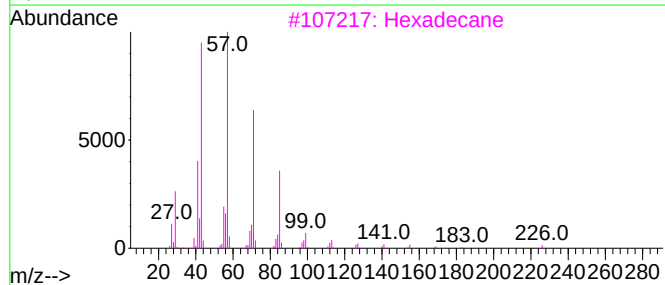
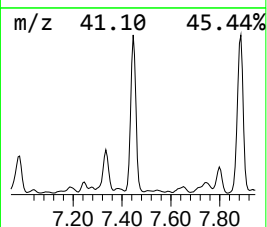
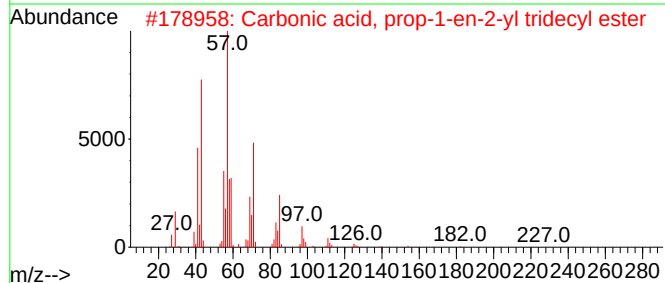
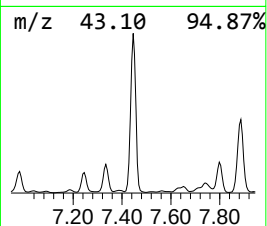
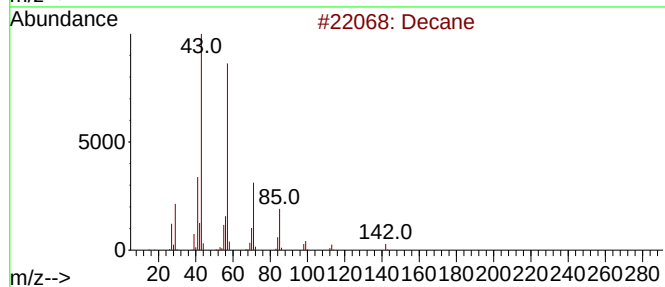
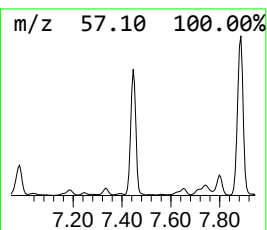
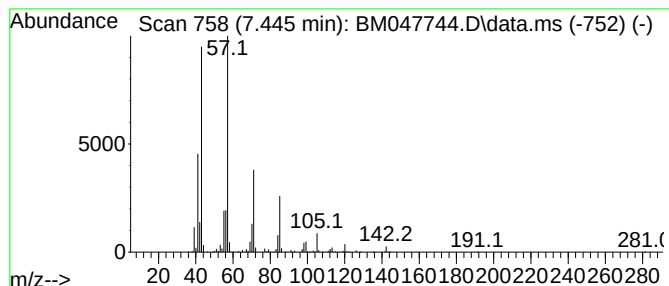
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.445	74.16 ng/ul	16660200	1,4-Dichlorobenzene-d4	7.810

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane	142	C10H22	000124-18-5	90
2		Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
3		Hexadecane	226	C16H34	000544-76-3	72
4		Tridecane	184	C13H28	000629-50-5	72
5		Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

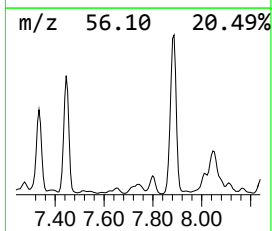
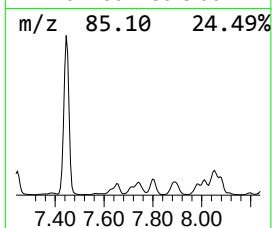
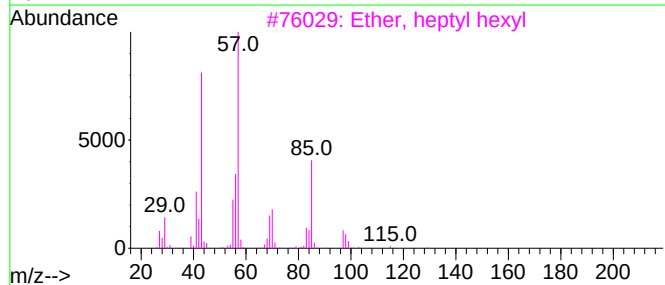
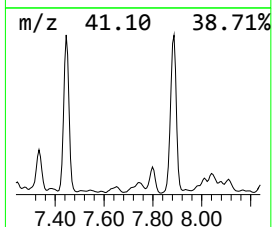
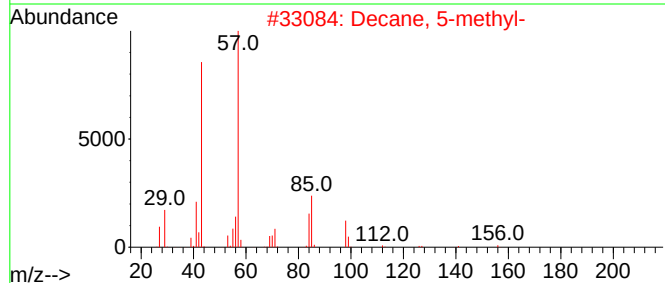
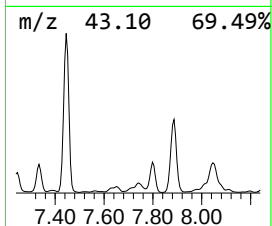
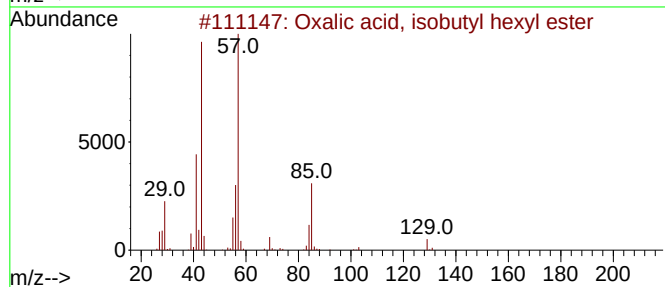
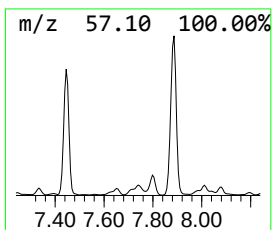
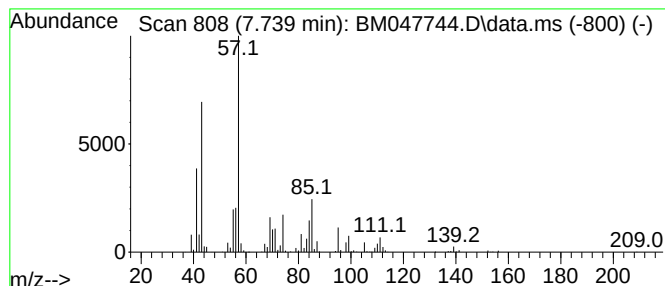
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Oxalic acid, isobutyl hexyl... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.739	8.96 ng/ul	2013420	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl hexyl ester	230	C12H22O4	1000309-37-1	50
2			Decane, 5-methyl-	156	C11H24	013151-35-4	49
3			Ether, heptyl hexyl	200	C13H28O	007289-40-9	47
4			Sulfurous acid, butyl hexyl ester	222	C10H22O3S	1000309-17-3	47
5			Heptane, 4-ethyl-	128	C9H20	002216-32-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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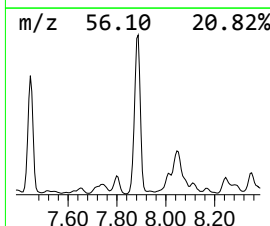
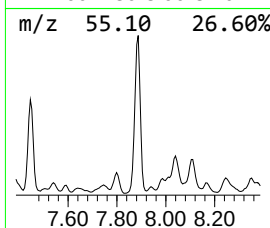
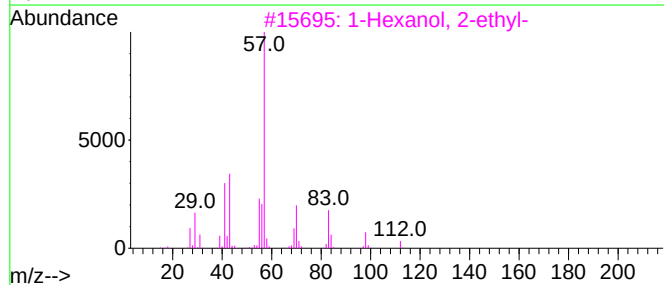
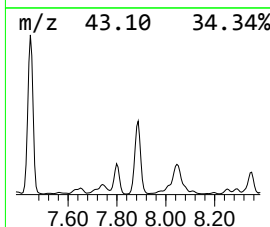
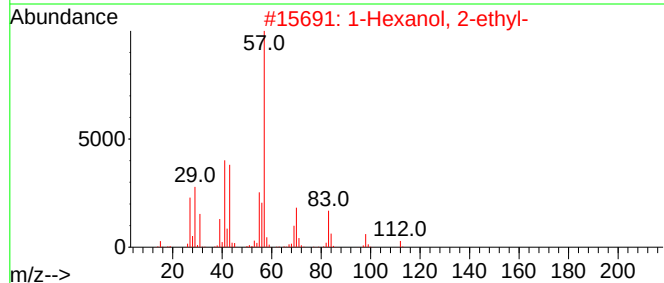
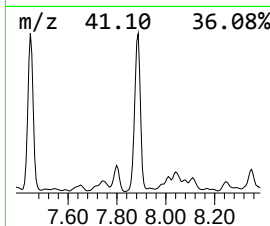
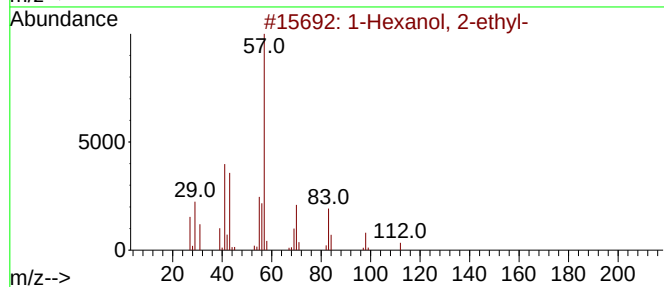
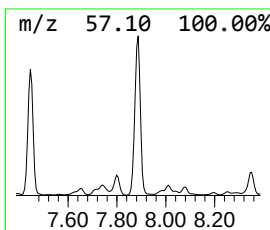
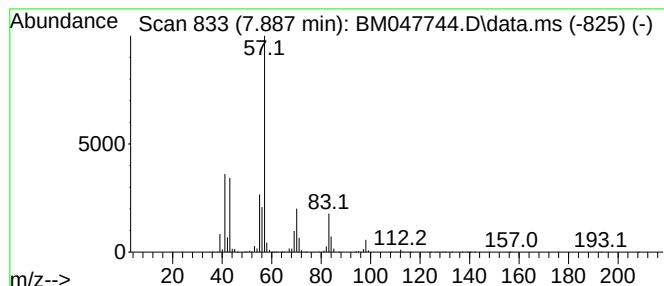
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	63.17 ng/ul	14190500	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Instrument :
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ClientSampleId :
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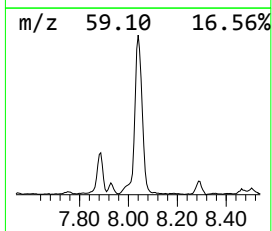
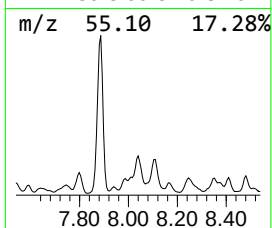
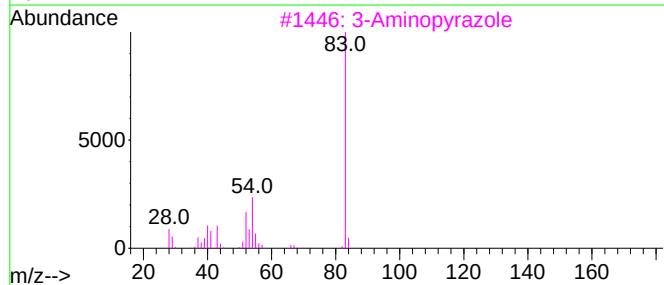
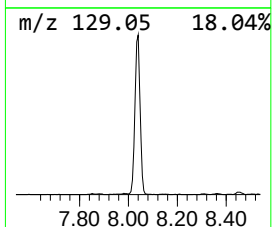
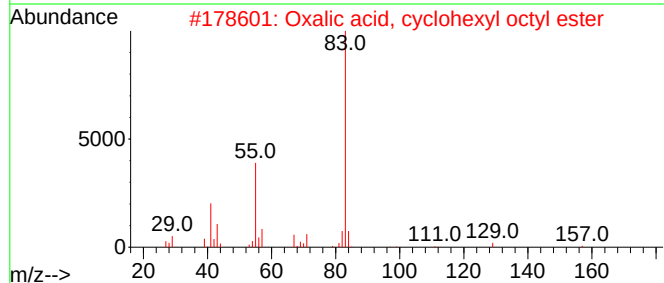
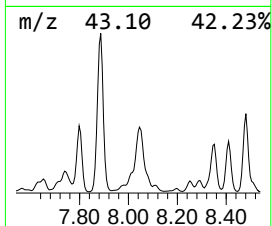
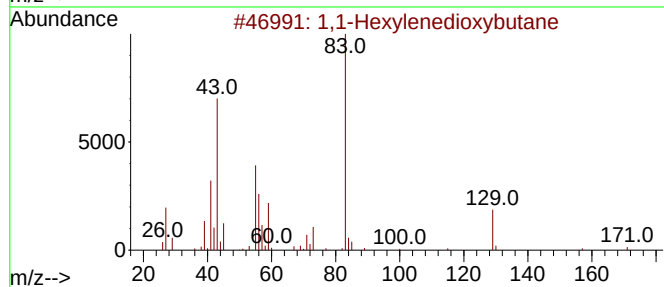
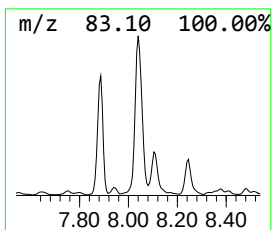
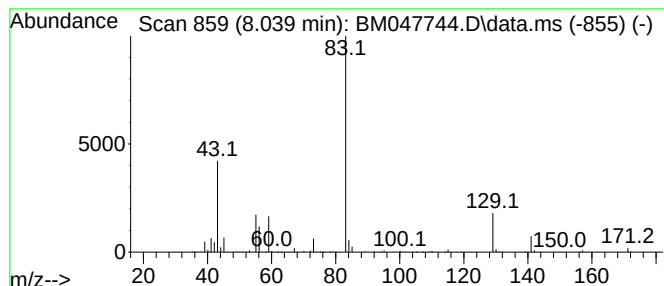
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1,1-Hexylenedioxybutane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.039	23.83 ng/ul	5353290	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	50
2			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	33
3			3-Aminopyrazole	83	C3H5N3	001820-80-0	9
4			4H-1,2,4-Triazole, 4-methyl-	83	C3H5N3	010570-40-8	9
5			1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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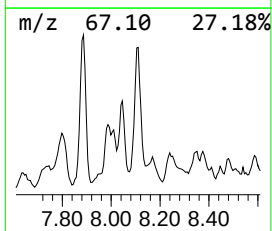
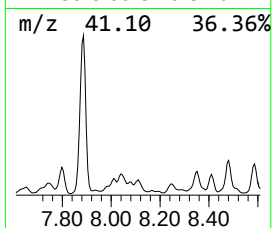
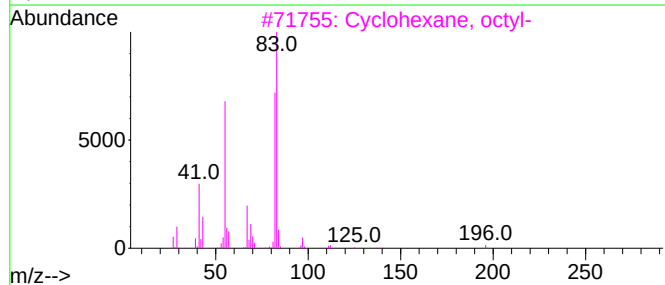
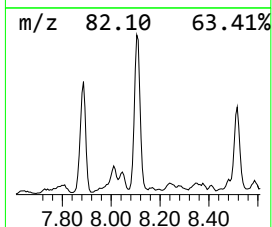
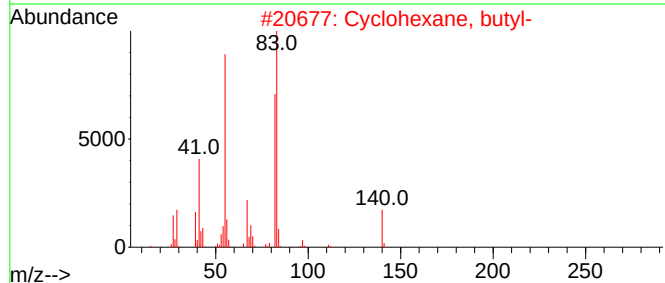
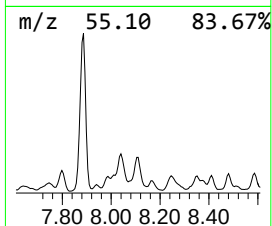
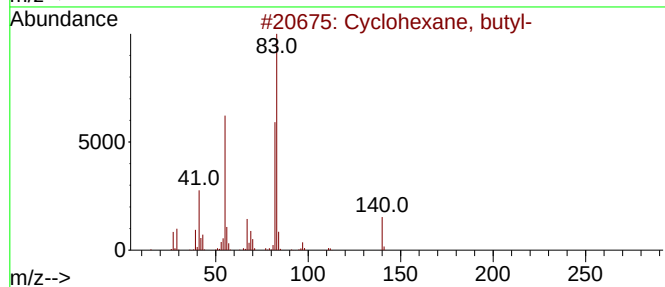
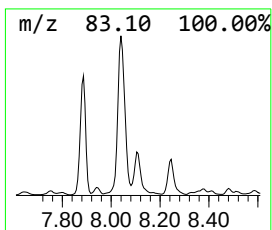
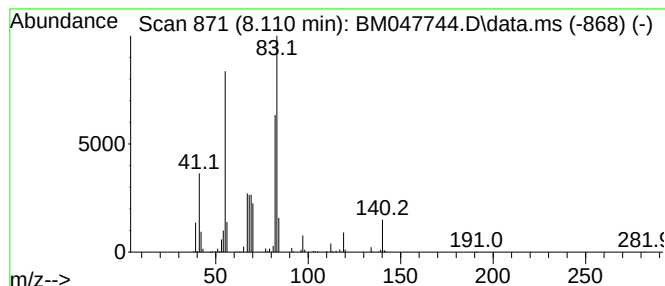
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Cyclic8.110 Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	6.25 ng/ul	1404640	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	68
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	64
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	59
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53
5			Cyclohexane, undecyl-	238	C17H34	054105-66-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Misc :
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ClientSampleId :
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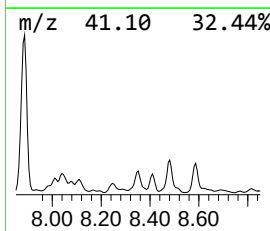
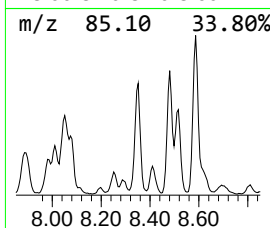
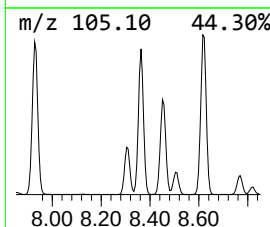
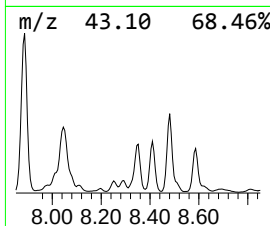
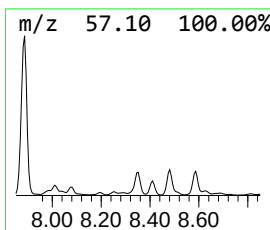
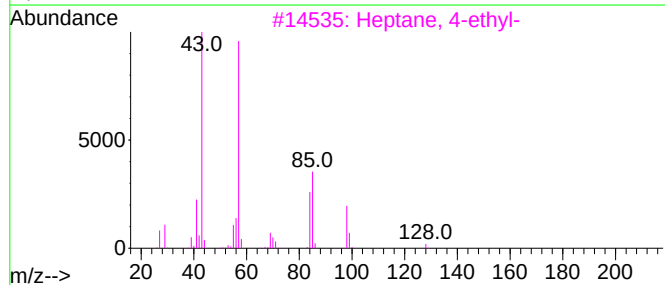
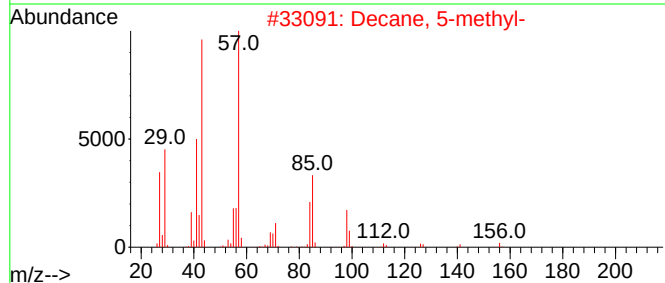
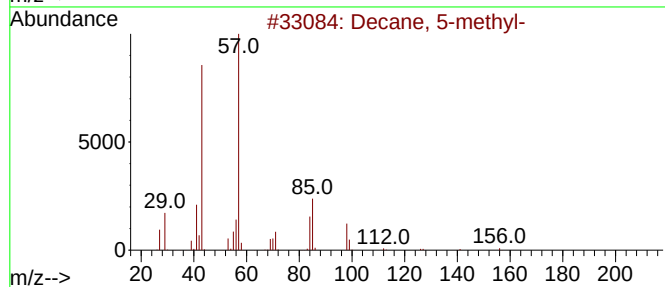
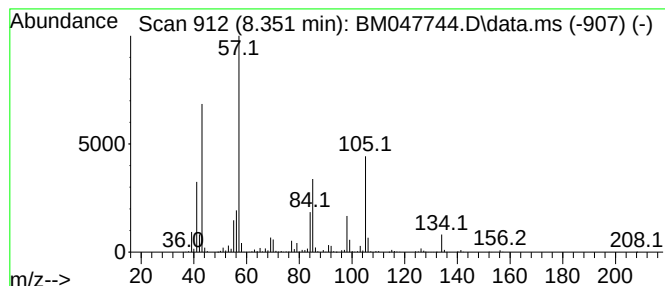
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.351	15.92 ng/ul	3576280	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	81
2			Decane, 5-methyl-	156	C11H24	013151-35-4	64
3			Heptane, 4-ethyl-	128	C9H20	002216-32-2	53
4			Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	50
5			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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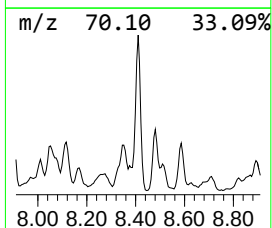
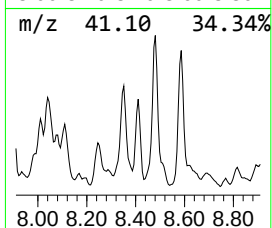
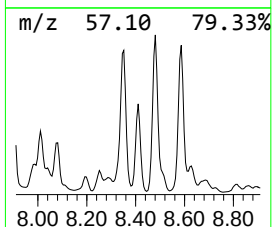
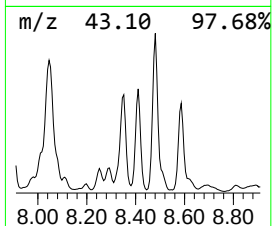
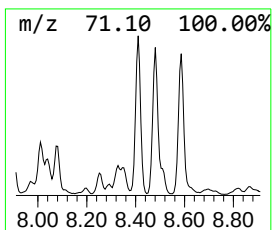
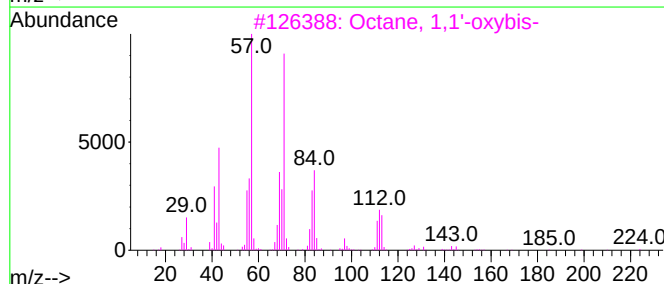
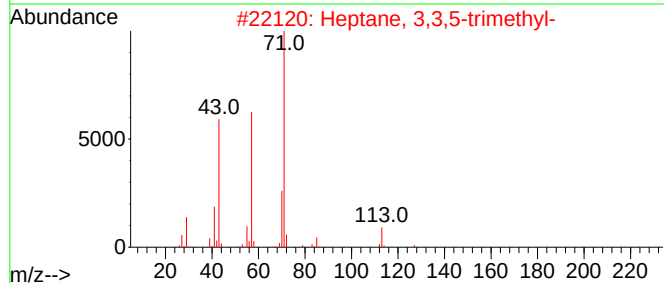
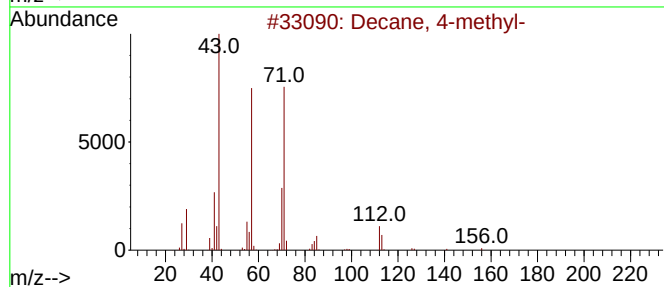
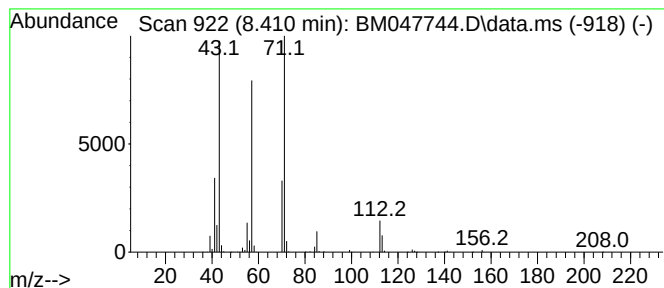
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.410	7.54 ng/ul	1692870	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	91
2			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72
3			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	64
4			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	64
5			Decane, 4-methyl-	156	C11H24	002847-72-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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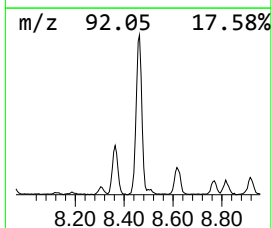
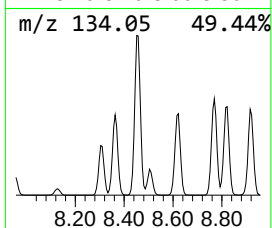
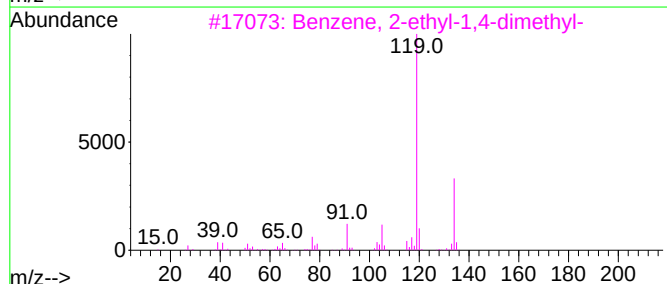
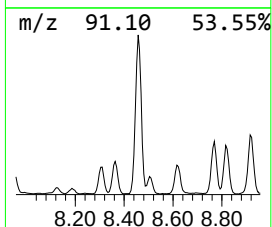
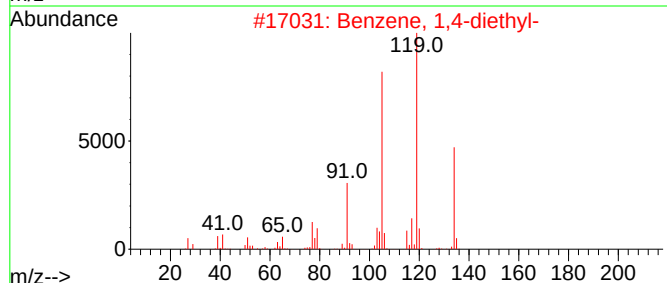
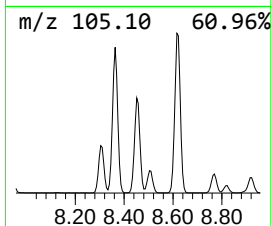
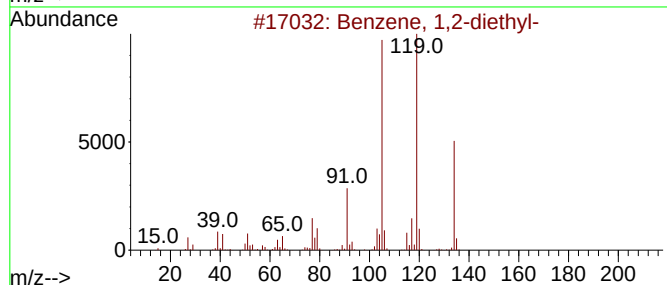
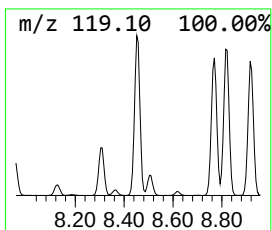
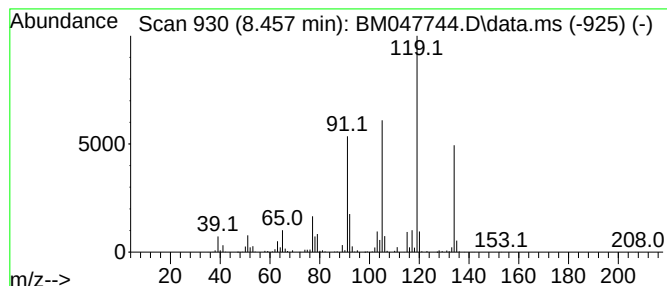
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 1,2-diethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.457	9.86 ng/ul	2214070	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
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Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

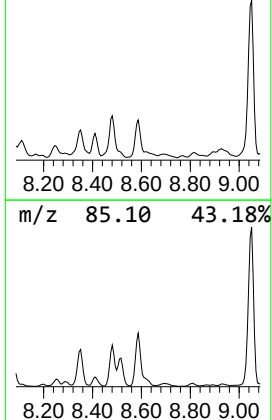
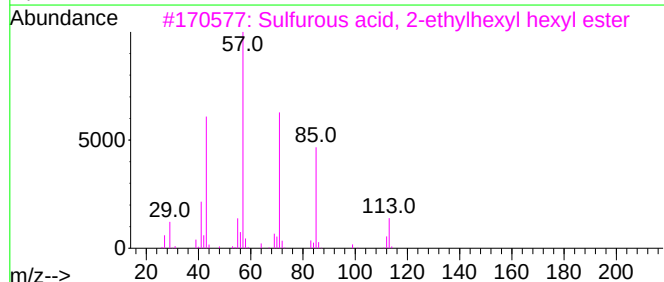
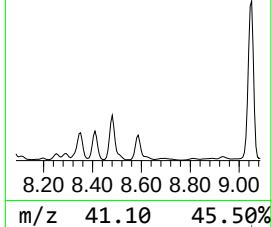
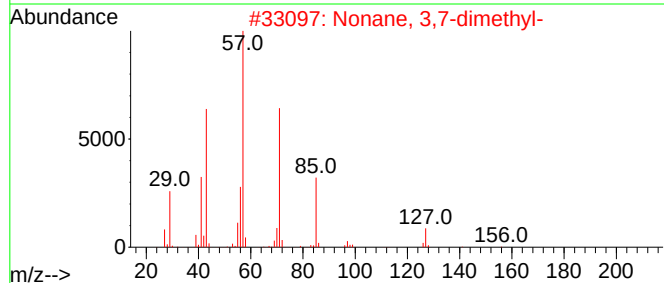
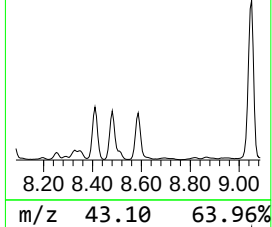
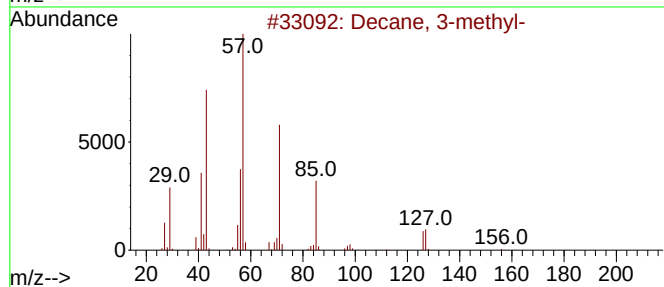
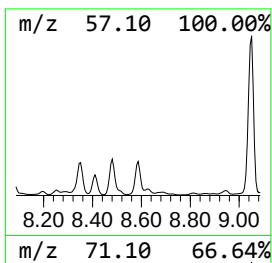
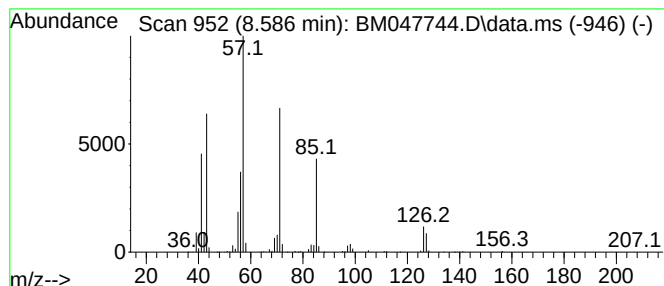
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.586	12.06 ng/ul	2708110	1,4-Dichlorobenzene-d4	7.810	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-	156	C11H24	013151-34-3	81
2	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	80
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	64
5	1-Iodoundecane	282	C11H23I	004282-44-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

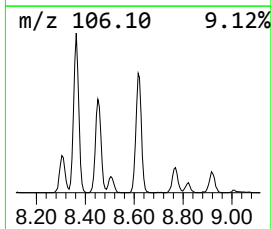
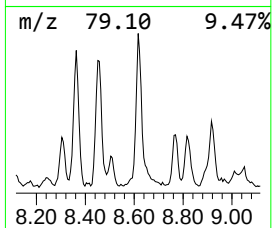
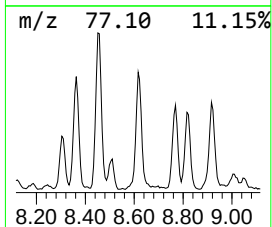
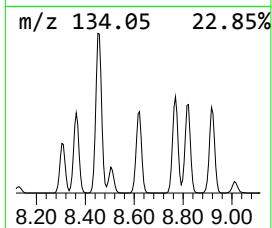
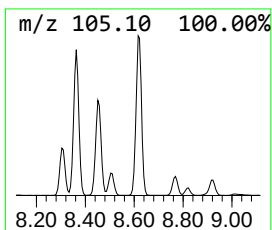
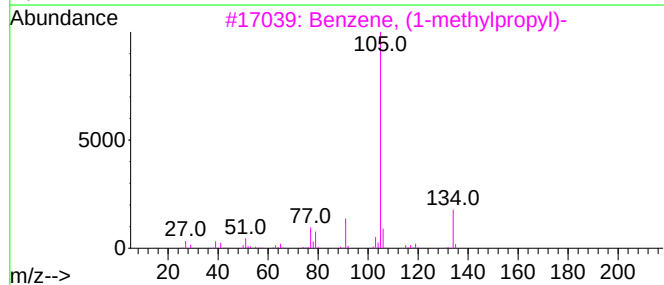
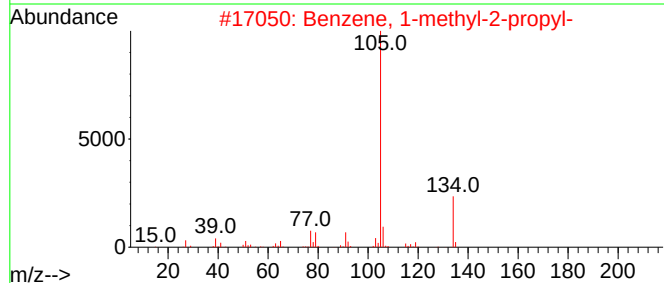
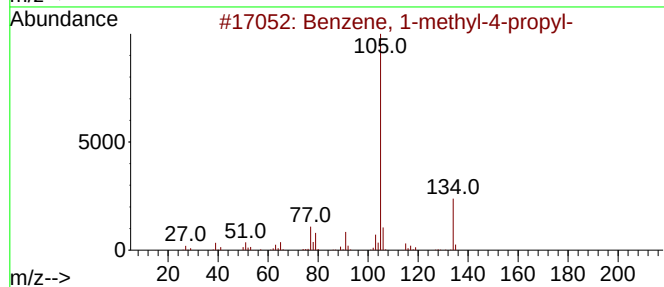
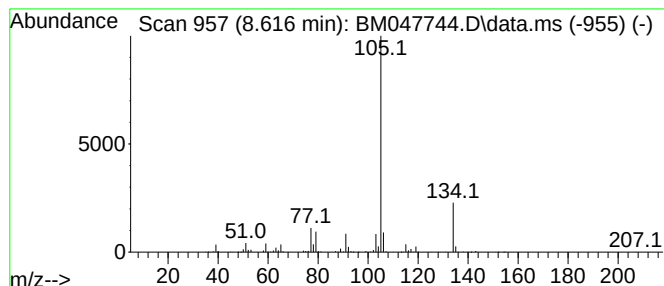
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Benzene, 1-methyl-4-propyl- Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.616	6.63 ng/ul	1489990	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	93
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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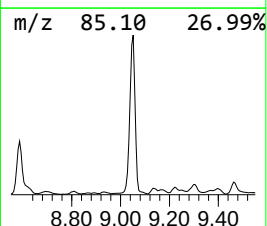
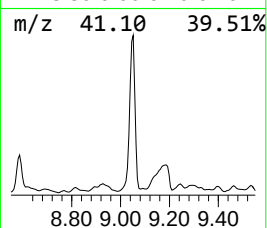
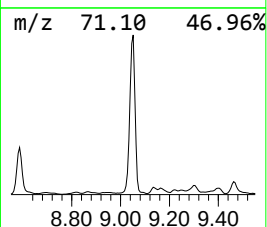
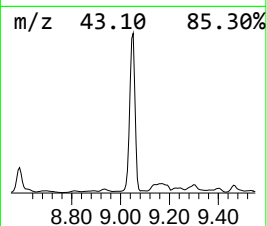
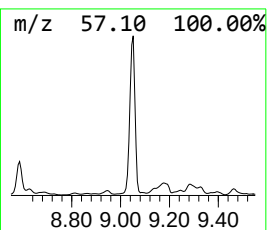
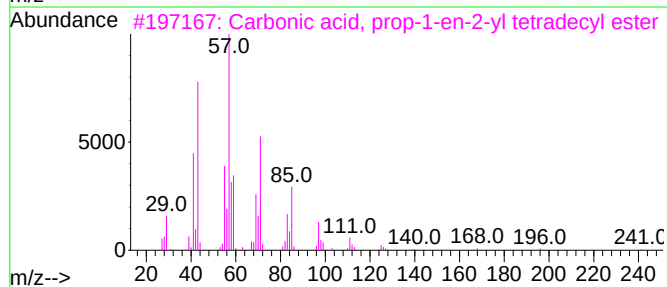
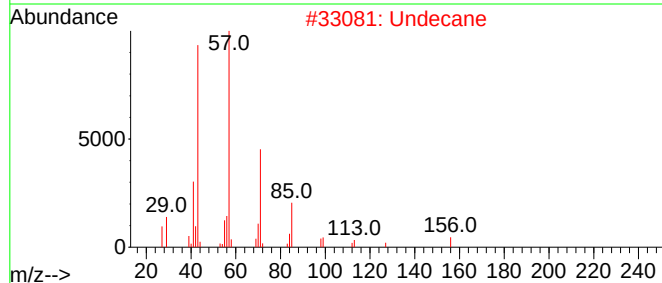
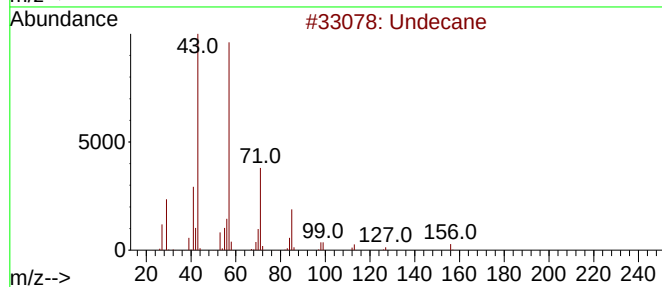
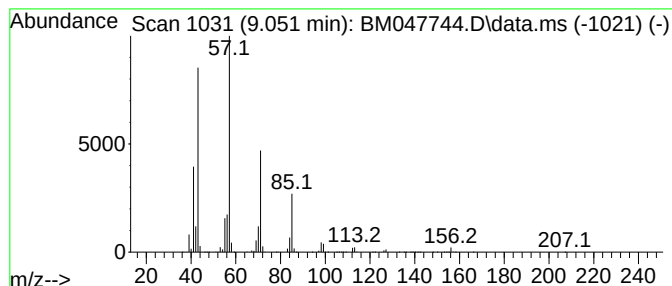
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.051	50.72 ng/ul	11393400	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Undecane	156	C11H24	001120-21-4	91
3			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
4			Tridecane	184	C13H28	000629-50-5	83
5			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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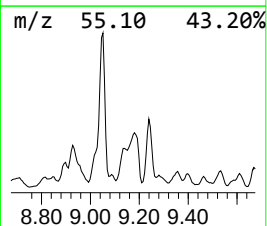
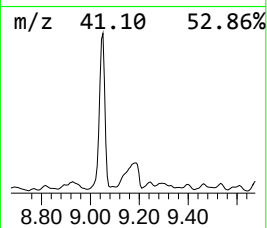
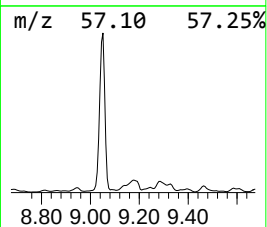
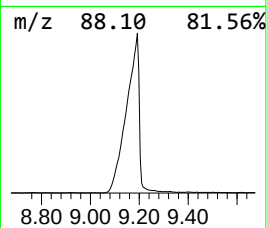
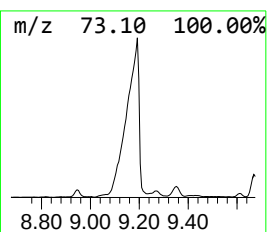
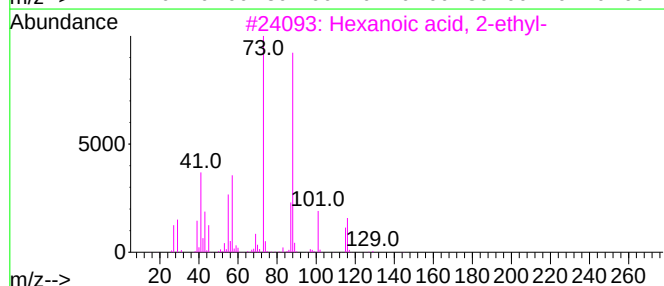
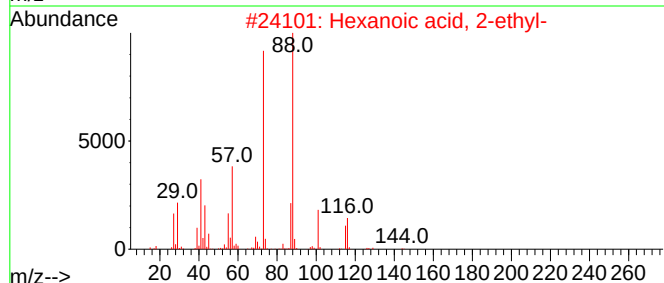
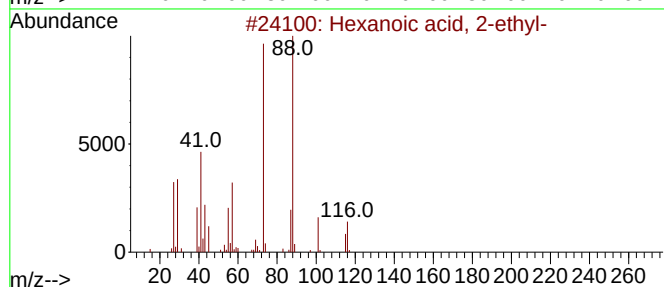
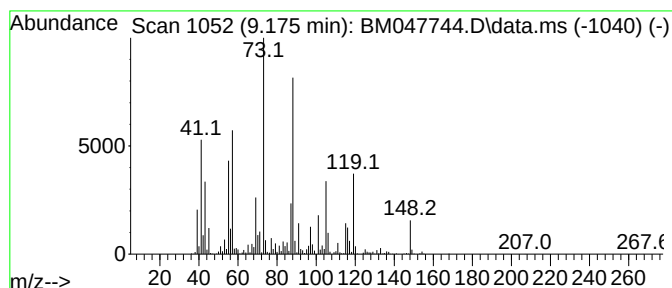
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.175	27.28 ng/ul	6128770	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	52
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	50
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	50
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	47
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
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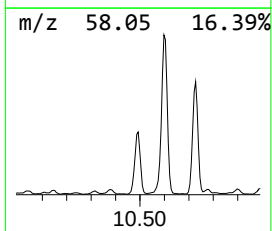
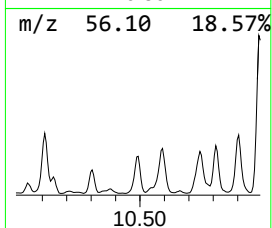
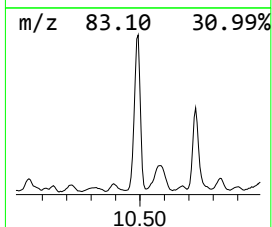
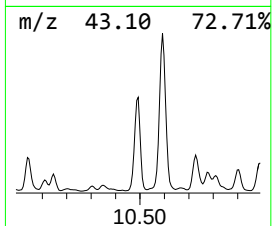
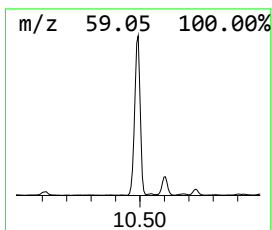
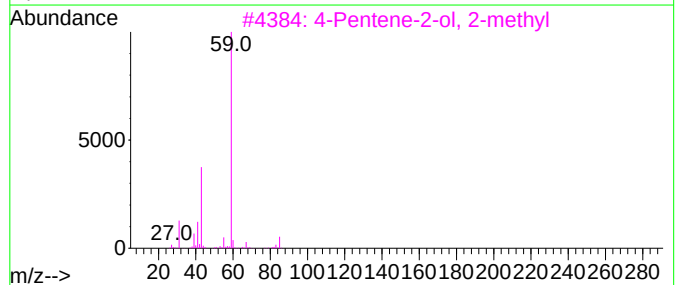
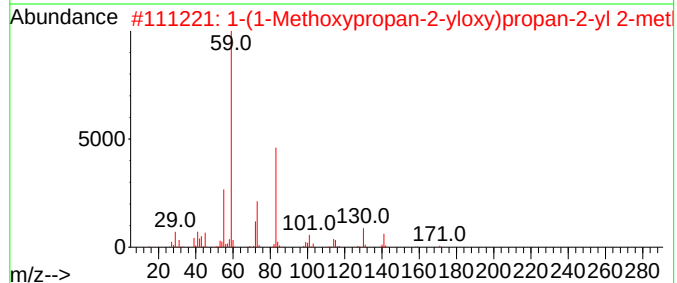
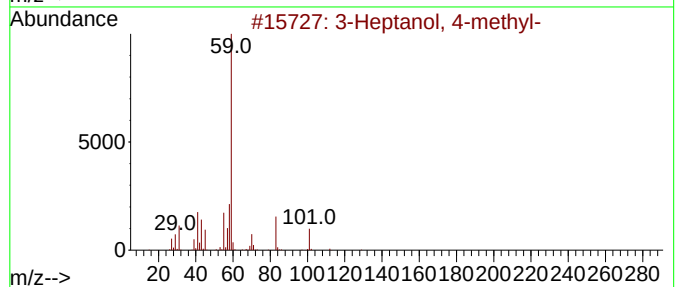
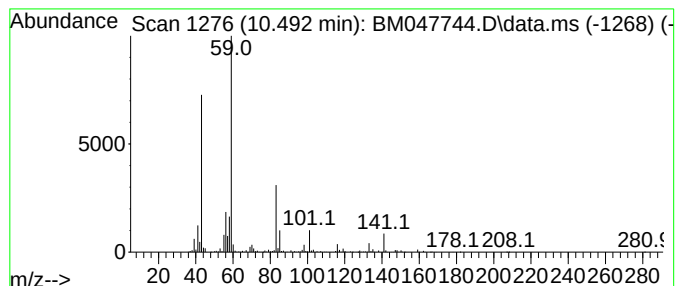
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 3-Heptanol, 4-methyl- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.492	9.11 ng/ul	4089110	Naphthalene-d8	10.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	59
2	1-(1-Methoxypropan-2-yloxy)propa...	230	C12H22O4	1000367-11-3	59
3	4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	47
4	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	47
5	Hexanamide	115	C6H13NO	000628-02-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
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Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

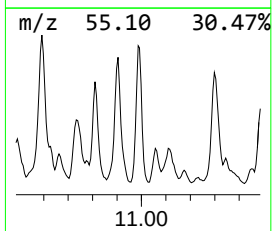
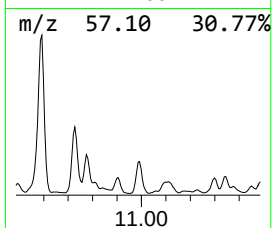
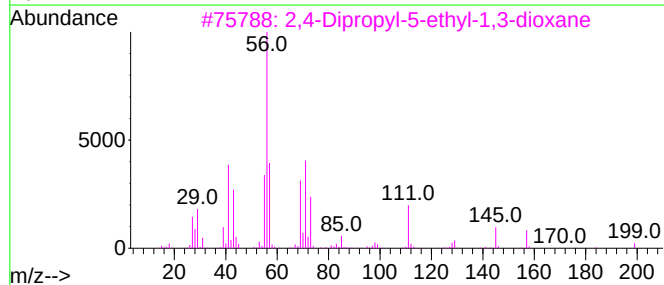
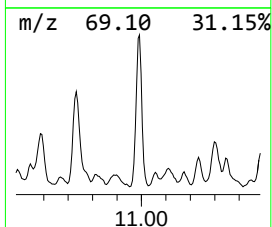
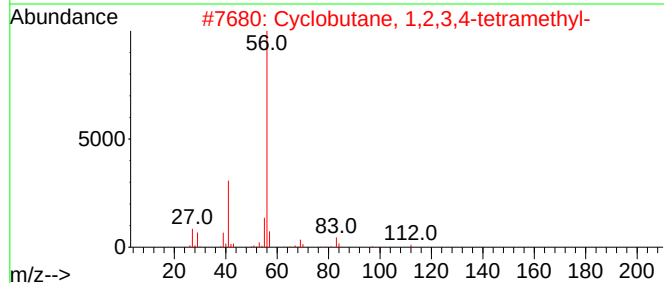
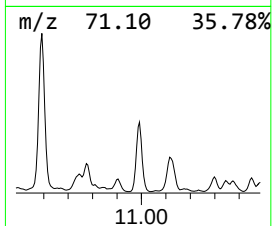
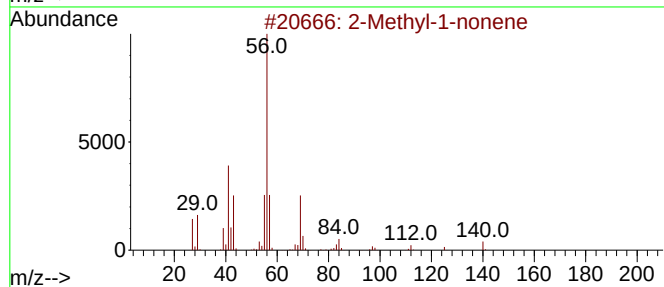
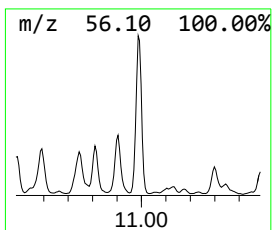
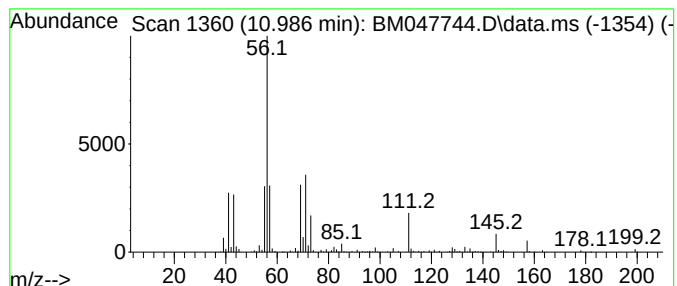
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.986	7.22 ng/ul	3242090	Naphthalene-d8	10.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-nonene	140	C10H20	002980-71-4	43
2	Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	43
3	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	37
4	Cyclohexylamine	99	C6H13N	000108-91-8	37
5	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

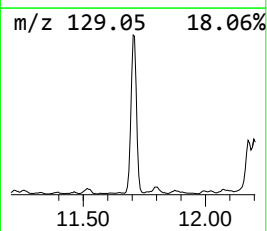
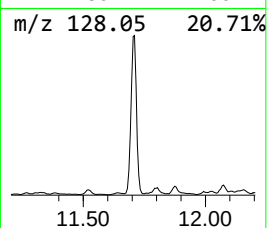
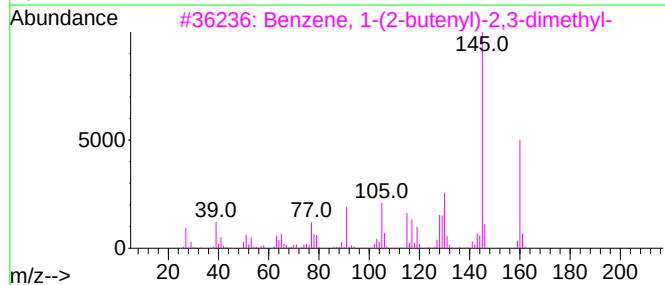
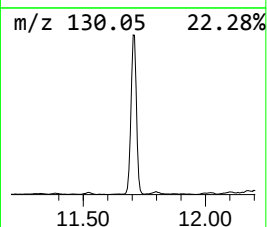
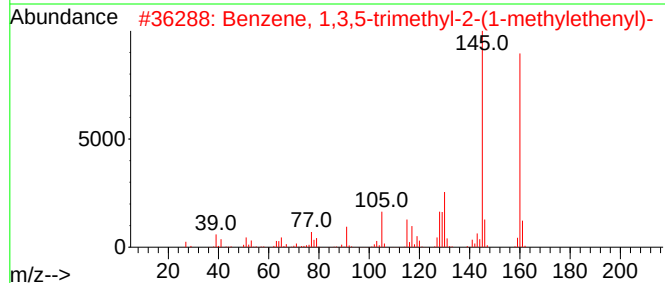
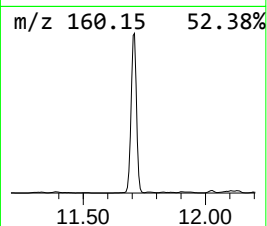
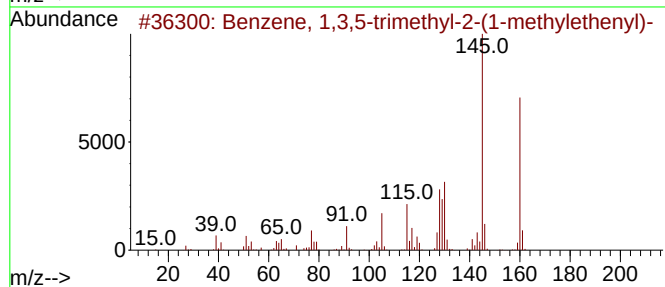
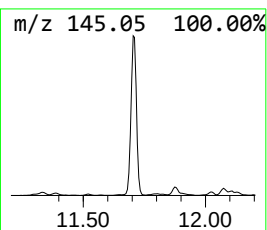
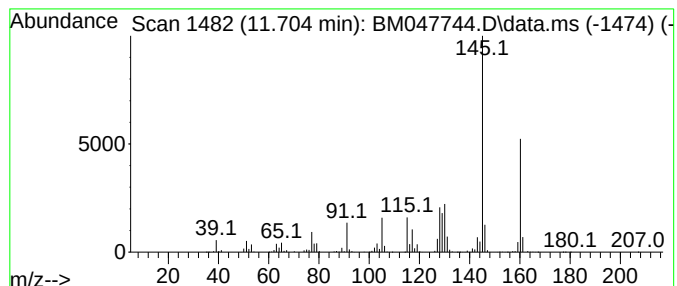
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	14.24 ng/ul	6395080	Naphthalene-d8	10.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	95
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	94
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
4			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	94
5			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

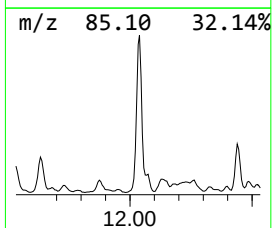
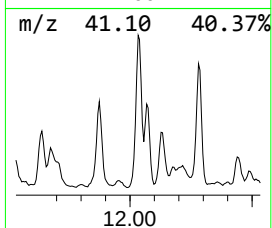
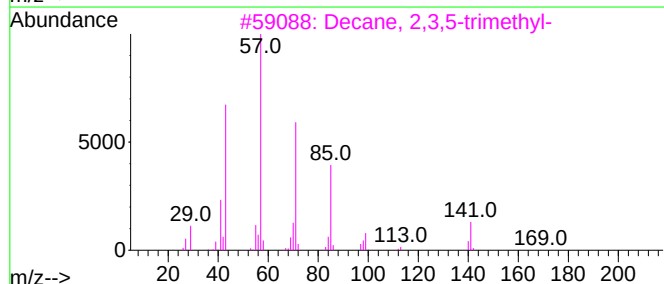
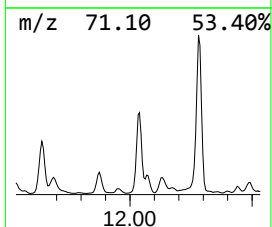
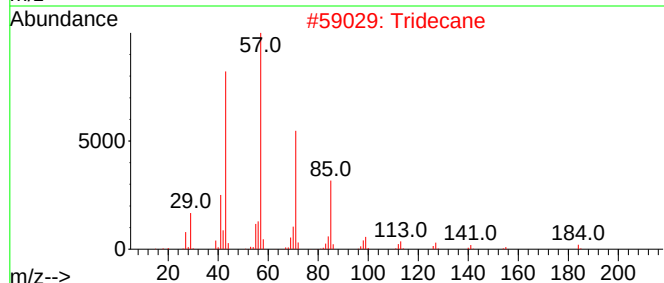
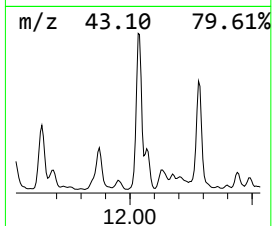
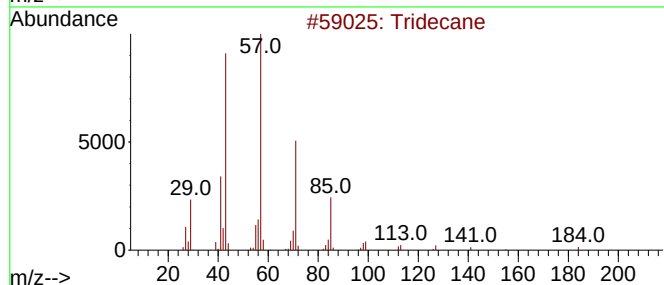
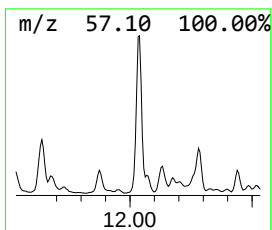
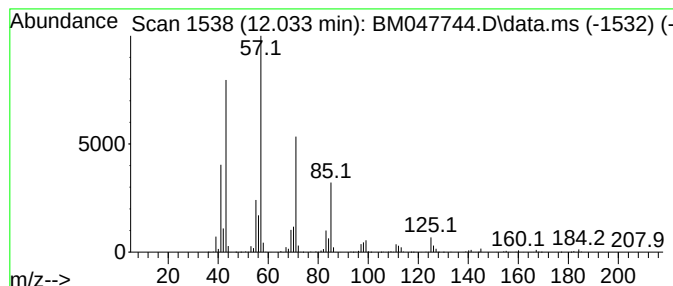
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.033	9.21 ng/ul	4137400	Naphthalene-d8	10.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	83
2	Tridecane	184	C13H28	000629-50-5	83
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	80
4	Tridecane	184	C13H28	000629-50-5	80
5	Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

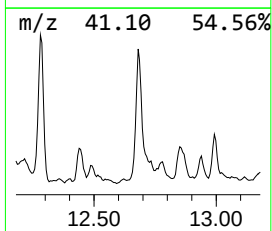
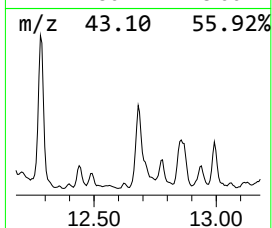
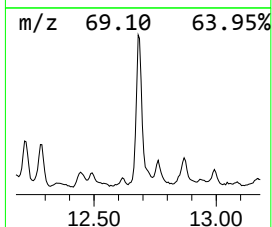
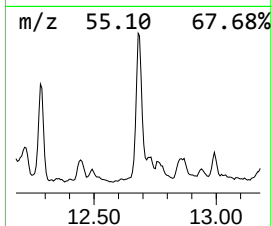
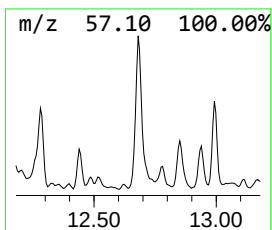
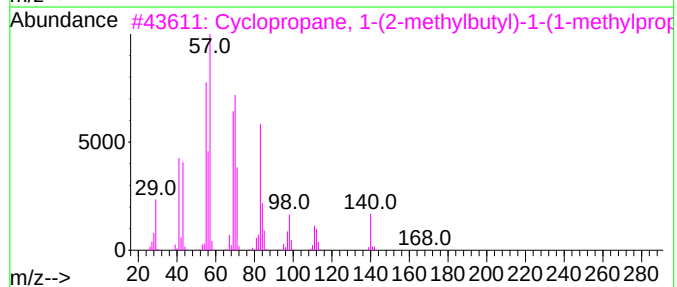
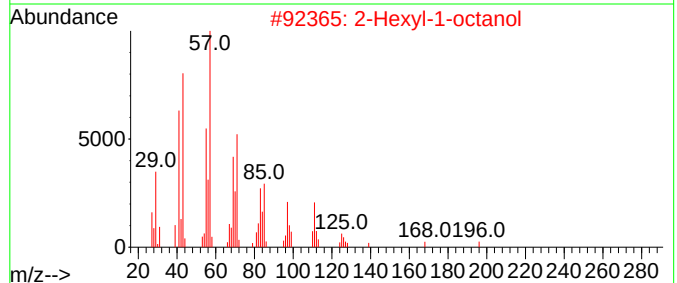
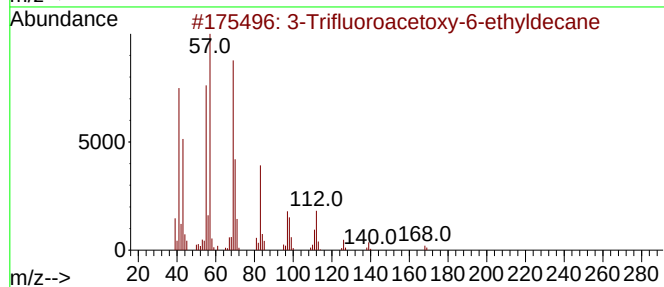
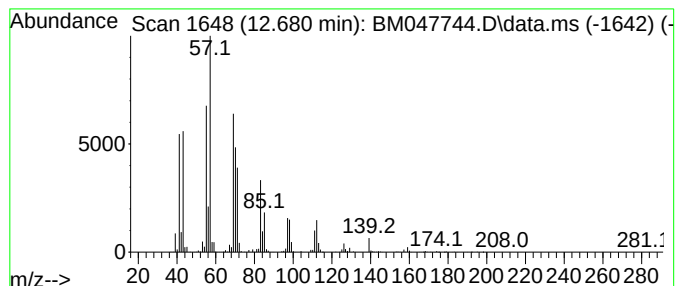
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 41 3-Trifluoroacetoxy-6-ethylde... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.680	20.25 ng/ul	3953870	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Trifluoroacetoxy-6-ethyldecane	282	C14H25F3O2	116436-59-0	58
2	2-Hexyl-1-octanol	214	C14H30O	019780-79-1	53
3	Cyclopropane, 1-(2-methylbutyl)-...	168	C12H24	064723-36-0	45
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	38
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

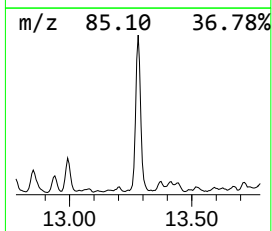
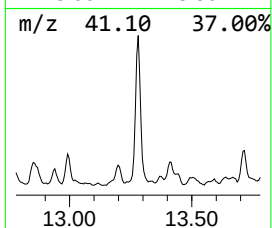
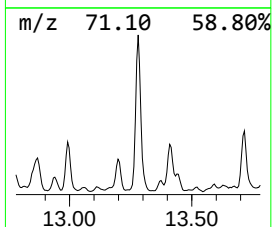
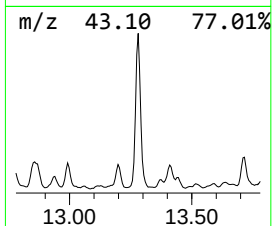
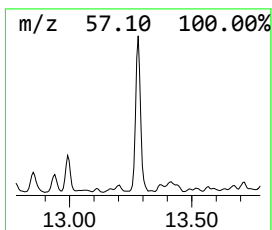
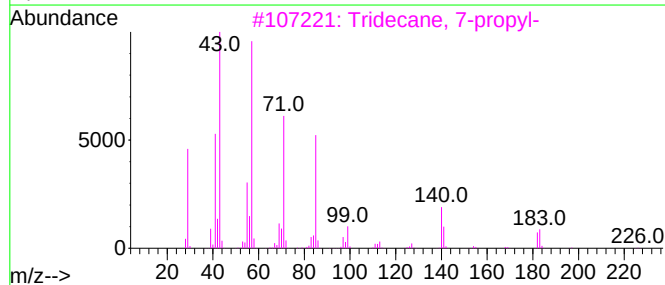
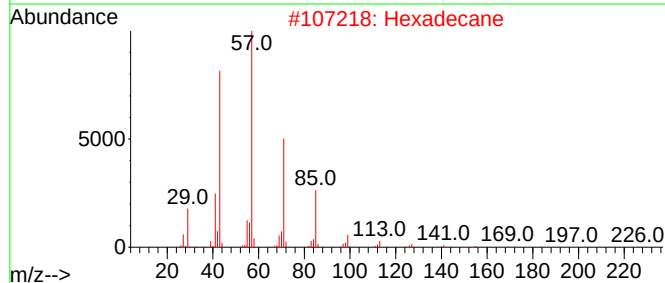
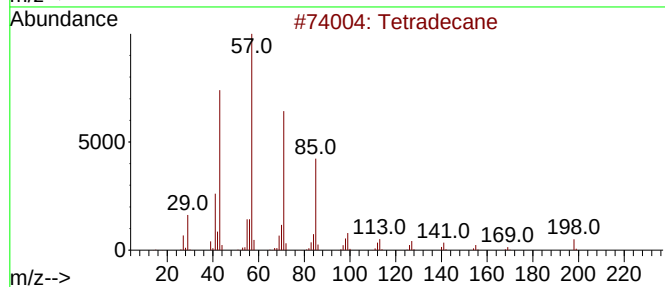
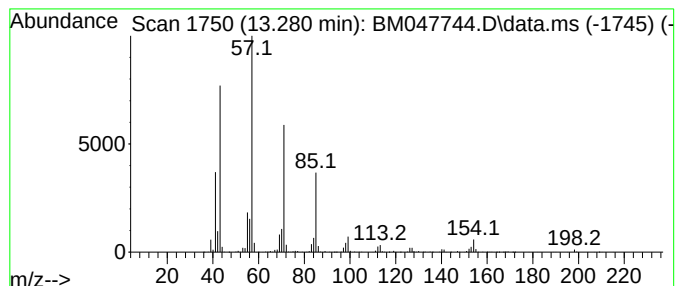
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.280	22.32 ng/ul	4357120	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	93
2	Hexadecane	226	C16H34	000544-76-3	87
3	Tridecane, 7-propyl-	226	C16H34	055045-09-5	86
4	Tridecane, 6-methyl-	198	C14H30	013287-21-3	86
5	Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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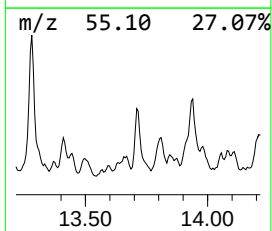
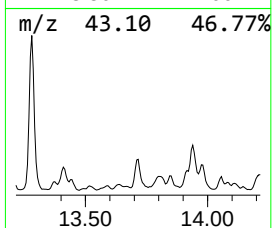
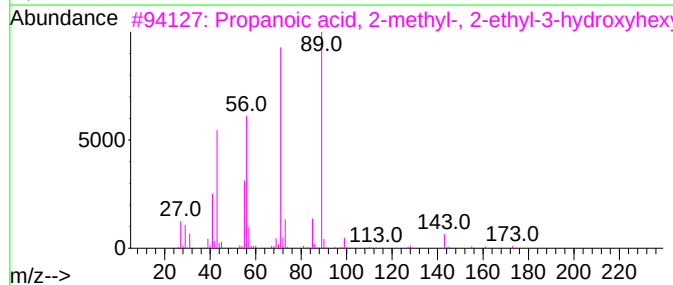
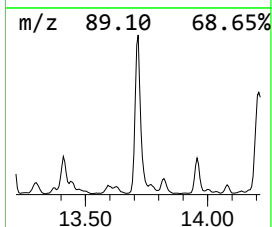
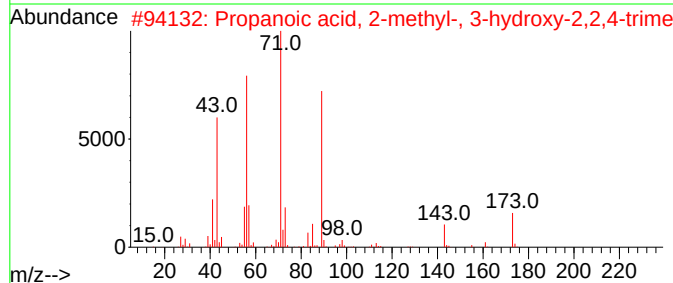
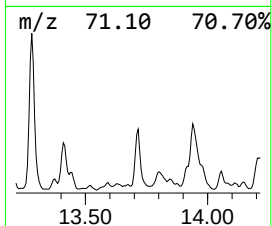
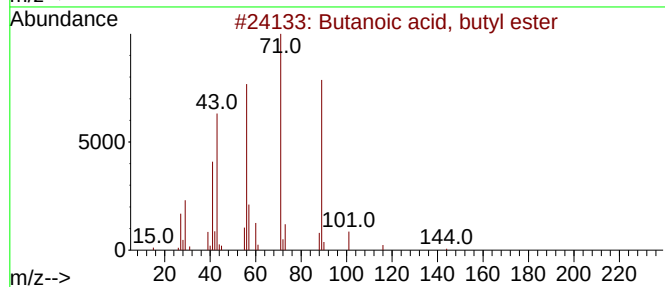
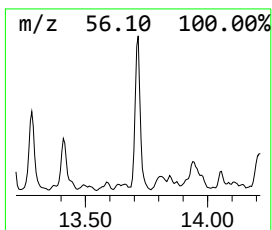
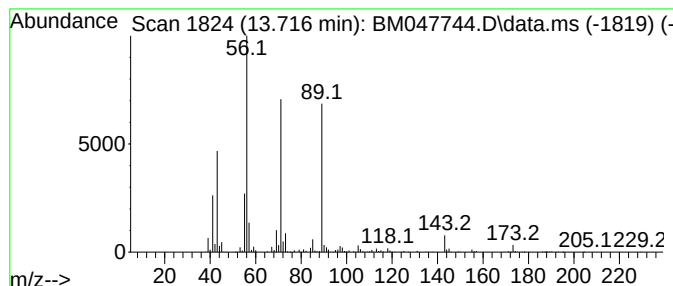
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 unknown-01 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	6.92 ng/ul	1350620	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	42
2			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	42
3			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	42
4			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	40
5			Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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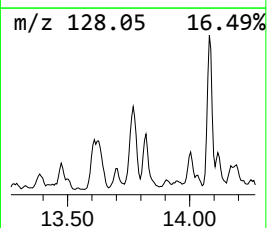
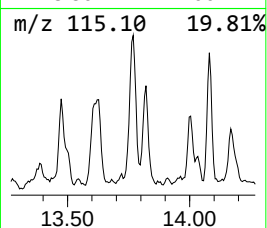
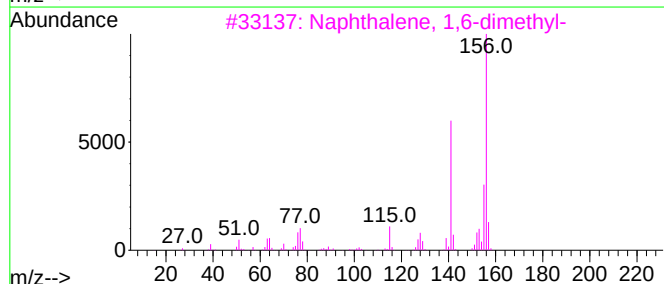
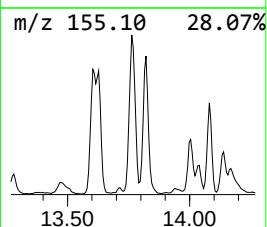
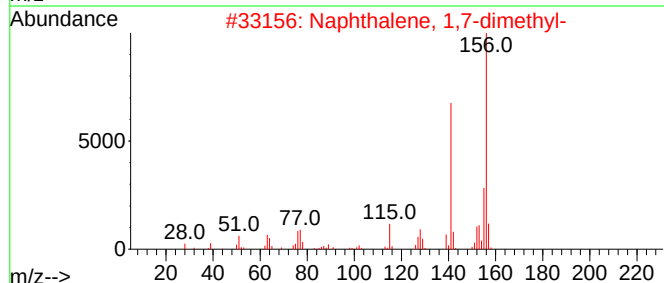
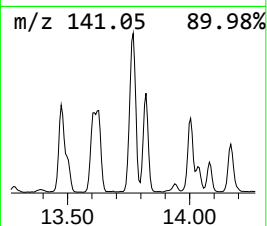
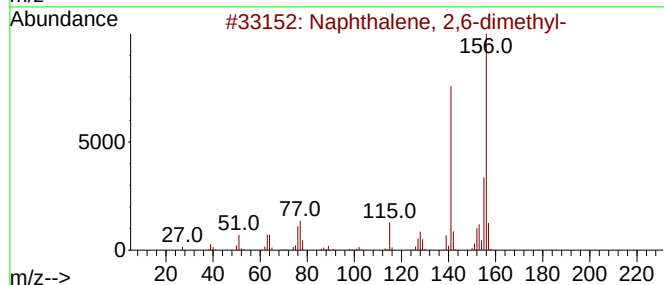
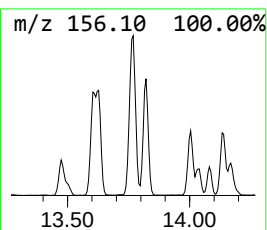
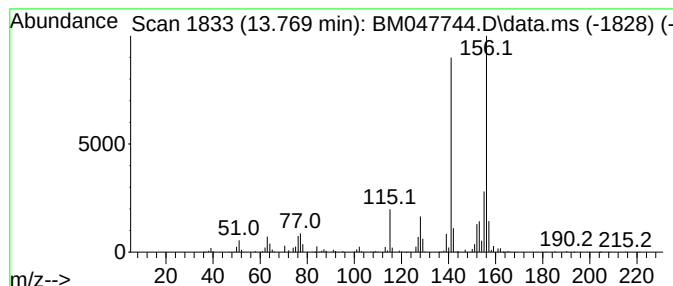
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 Naphthalene, 2,6-dimethyl- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	8.64 ng/ul	1687830	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

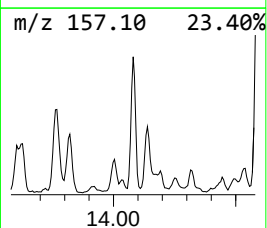
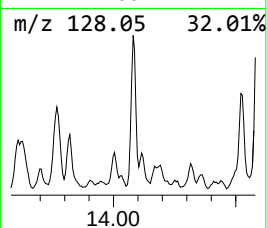
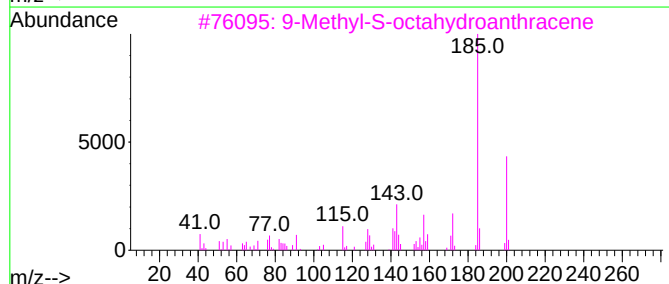
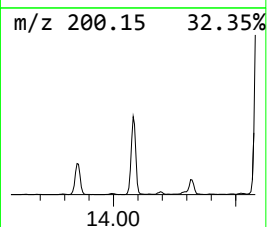
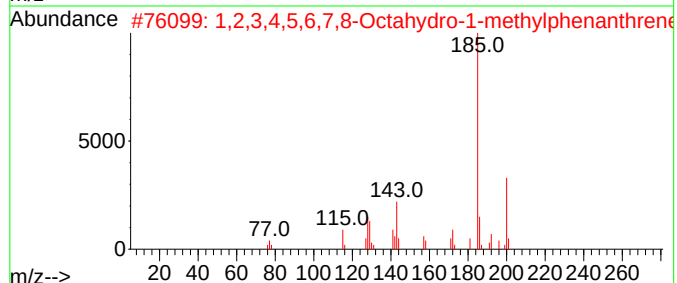
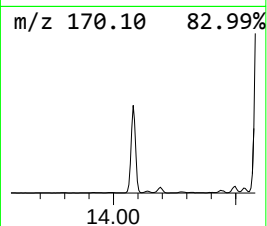
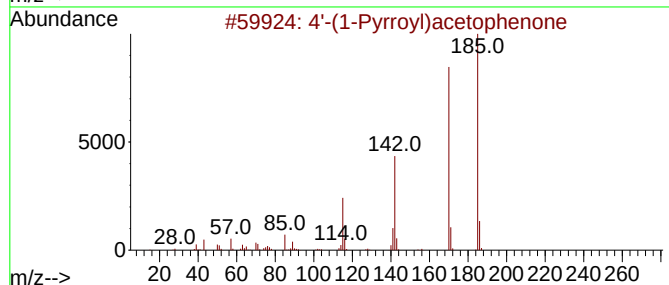
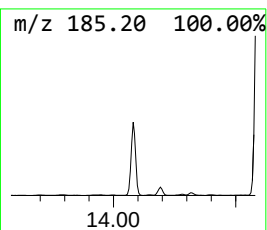
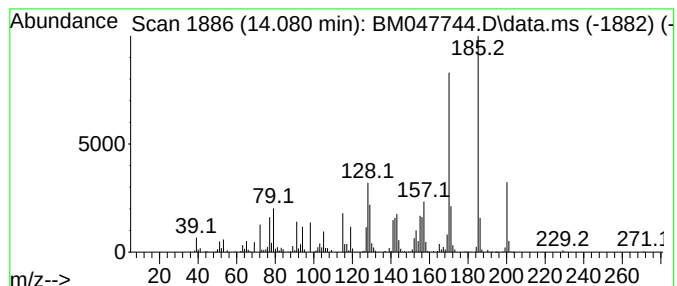
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 unknown-02 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	10.13 ng/ul	1978240	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4'-(1-Pyrroyl)acetophenone		185	C12H11NO	022106-37-2	49	
2	1,2,3,4,5,6,7,8-Octahydro-1-meth...		200	C15H20	1000080-19-4	49	
3	9-Methyl-S-octahydroanthracene		200	C15H20	004948-51-0	46	
4	Ethanone, 1-(5-methyl-1-phenyl-1...		200	C12H12N2O	006123-63-3	30	
5	Naphthalene, 1,4,6-trimethyl-		170	C13H14	002131-42-2	27	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

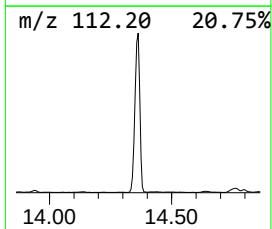
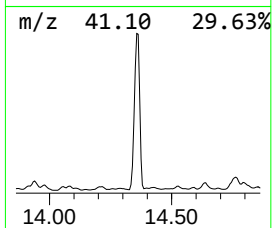
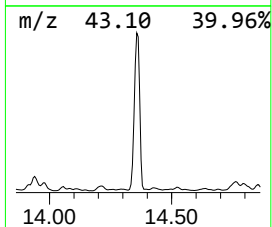
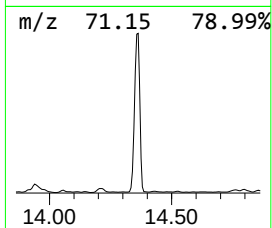
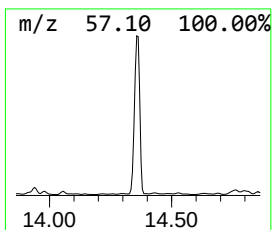
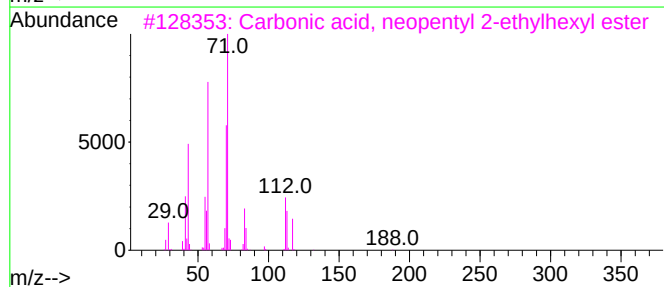
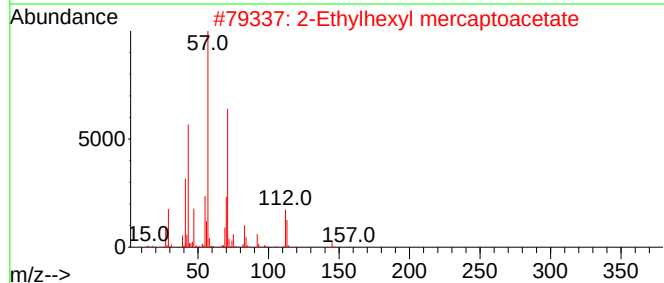
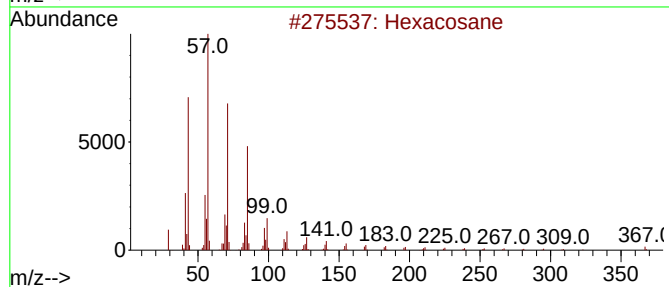
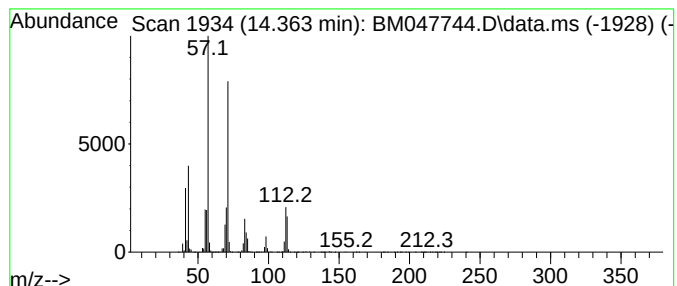
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.363	132.78 ng/ul	25924800	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	72
2	2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	72
3	Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	72
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
5	1-Decene, 4-methyl-	154	C11H22	013151-29-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

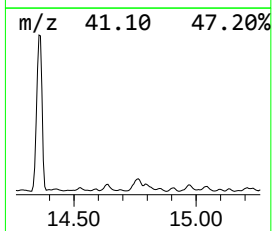
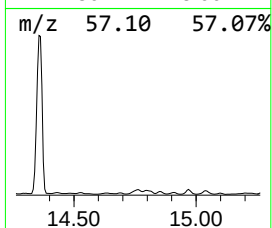
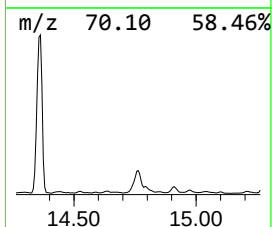
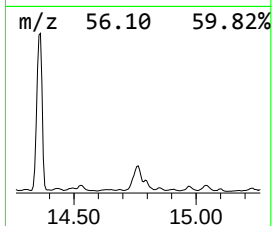
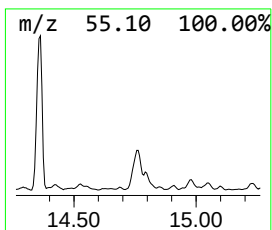
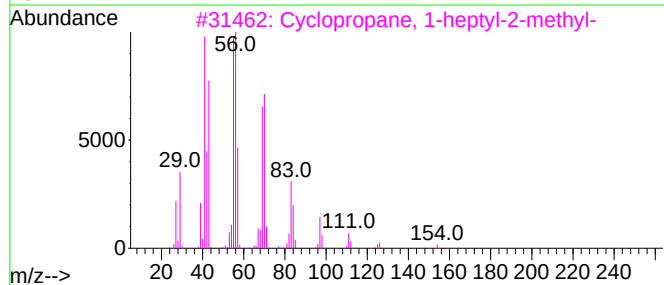
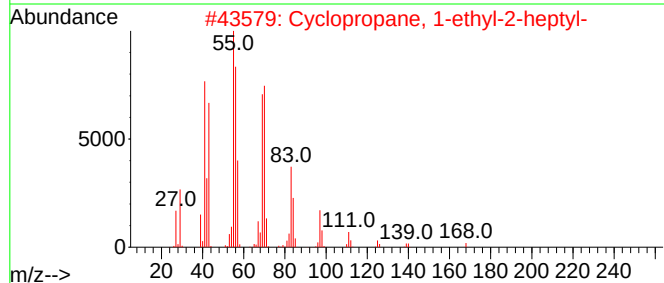
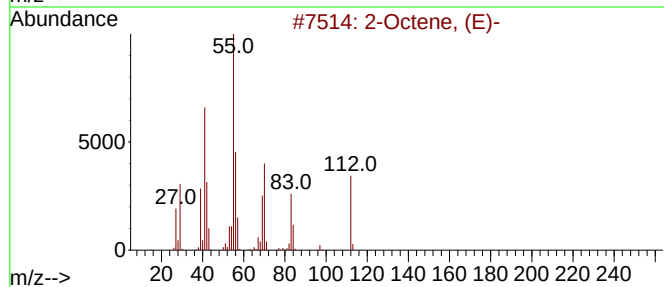
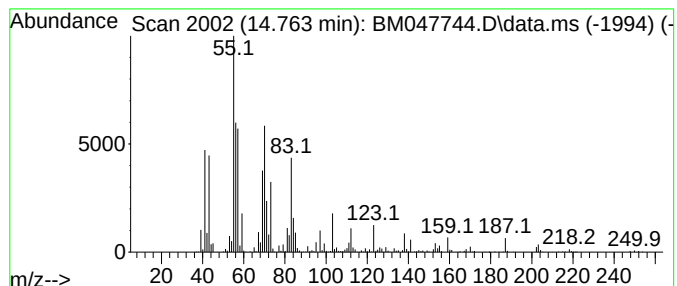
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.763	17.46 ng/ul	3409310	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Octene, (E)-		112	C8H16	013389-42-9	58
2	Cyclopropane, 1-ethyl-2-heptyl-		168	C12H24	074663-86-8	53
3	Cyclopropane, 1-heptyl-2-methyl-		154	C11H22	074663-91-5	46
4	Cyclopentane, 1,2,3-trimethyl-		112	C8H16	002815-57-8	38
5	1-Undecene, 9-methyl-		168	C12H24	074630-41-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

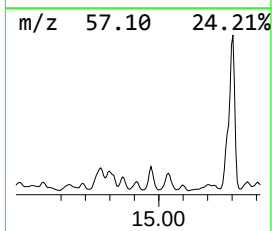
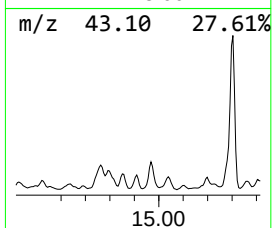
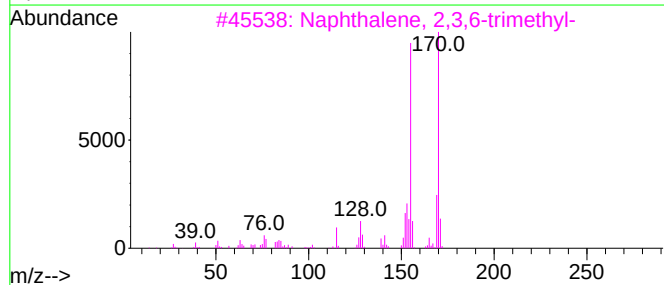
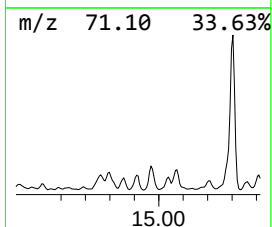
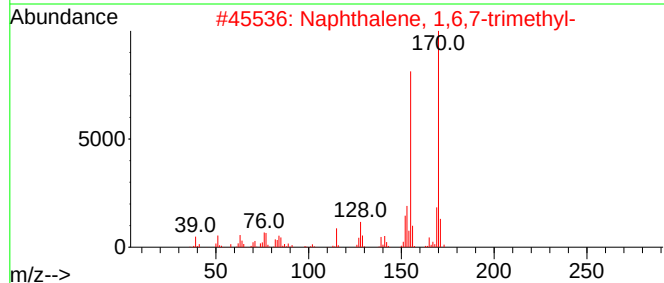
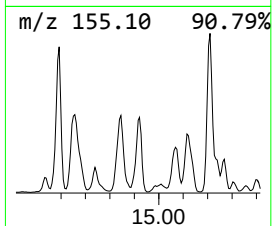
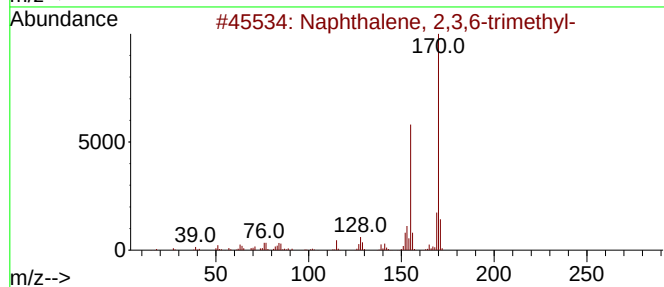
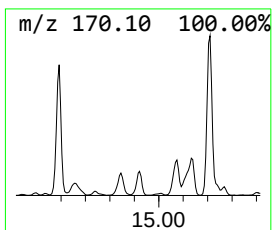
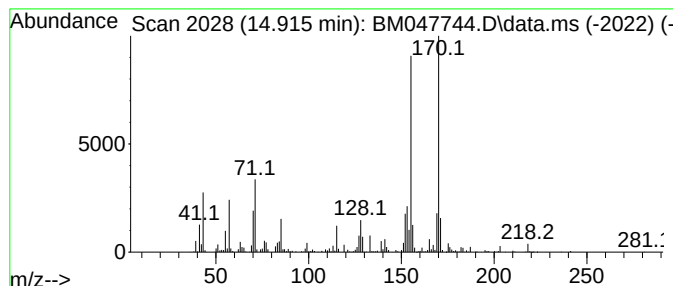
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 Naphthalene, 2,3,6-trimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.916	8.06 ng/ul	1574440	Acenaphthene-d10	14.445

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

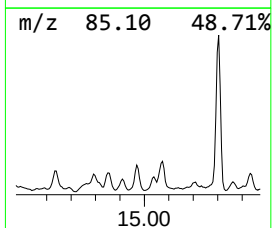
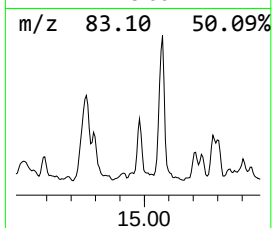
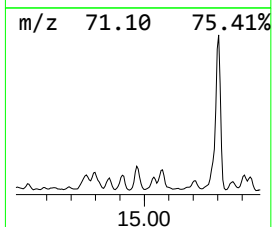
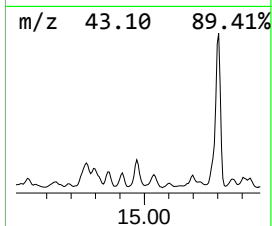
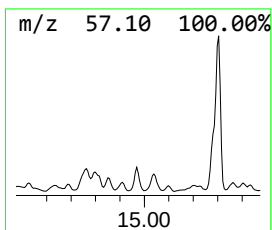
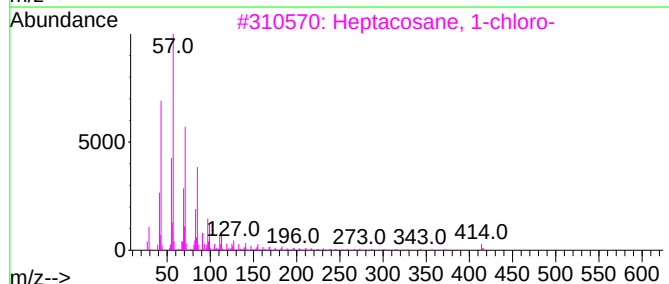
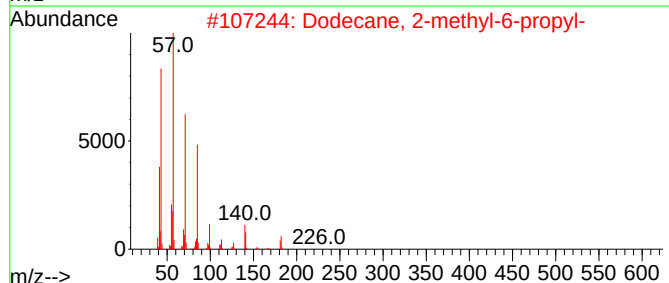
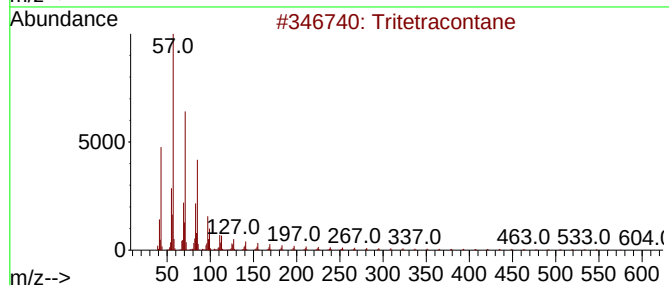
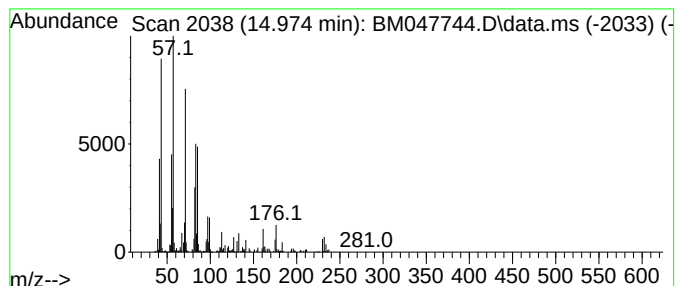
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.974	7.20 ng/ul	1405180	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tritetracontane	605	C43H88	007098-21-7	64
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	64
3	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	64
4	Tetratetracontane	619	C44H90	007098-22-8	62
5	Heneicosane	296	C21H44	000629-94-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

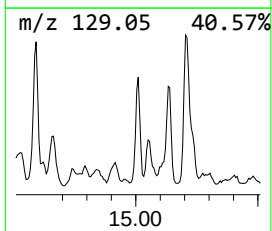
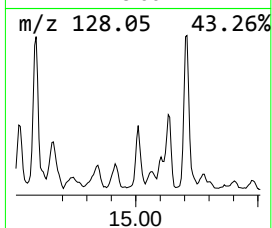
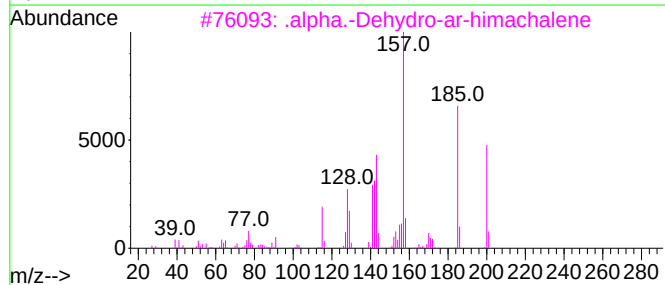
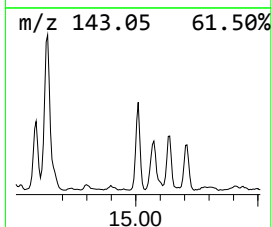
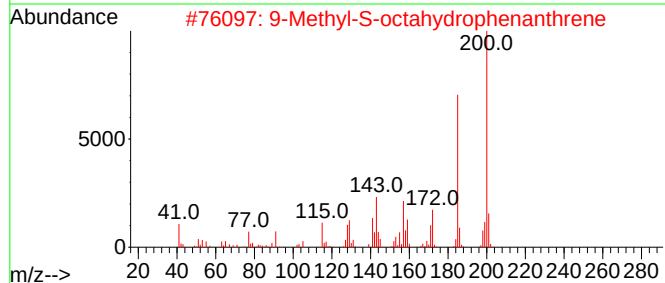
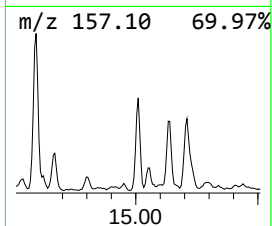
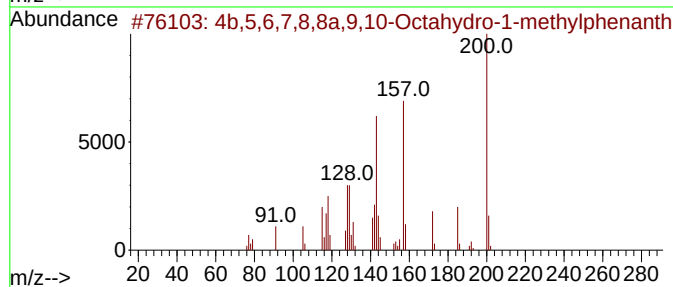
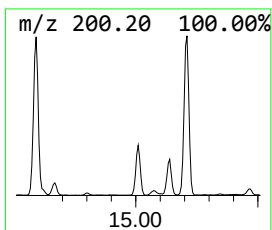
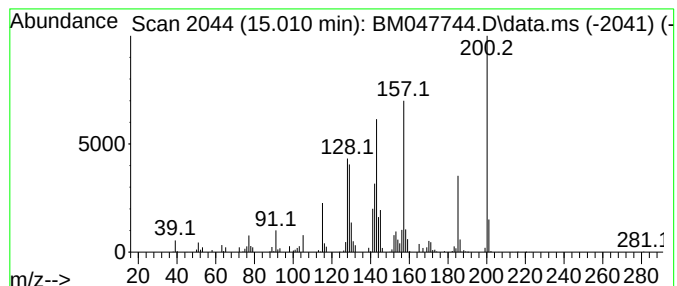
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 50 4b,5,6,7,8,8a,9,10-Octahydr... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.010	6.80 ng/ul	1328330	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4b,5,6,7,8,8a,9,10-Octahydro-1-m...		200	C15H20	1000080-19-3	70
2	9-Methyl-S-octahydrophenanthrene		200	C15H20	023861-81-6	62
3	.alpha.-Dehydro-ar-himachalene		200	C15H20	078204-62-3	59
4	.alpha.-Calacorene		200	C15H20	021391-99-1	55
5	4-Phenyl-2-(2-oxopropyl)furan		200	C13H12O2	1000426-70-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

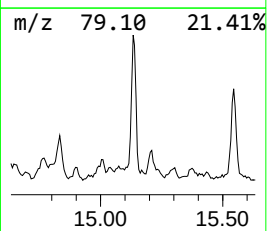
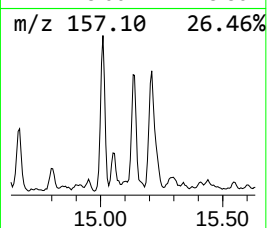
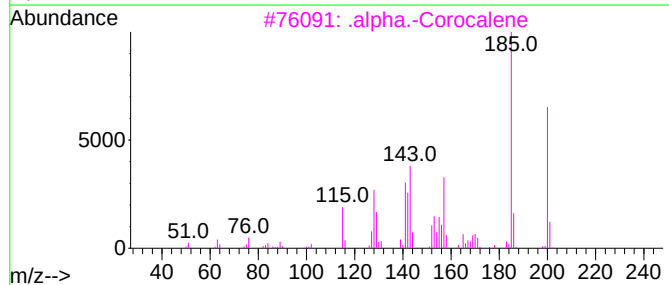
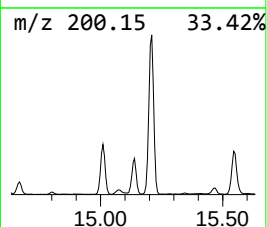
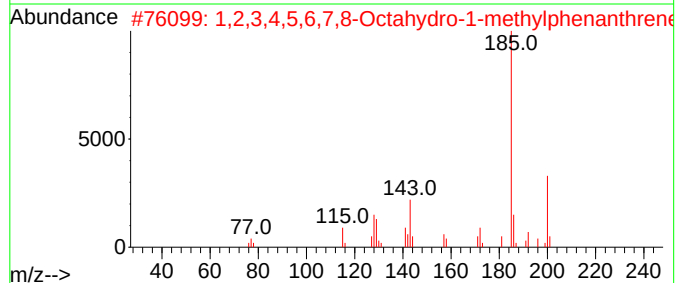
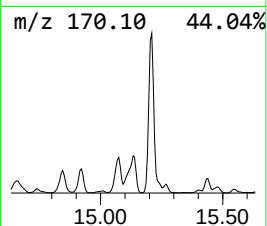
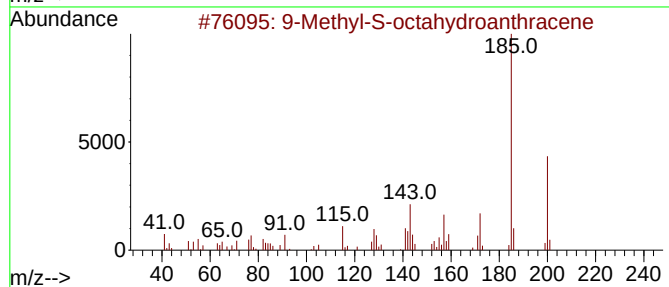
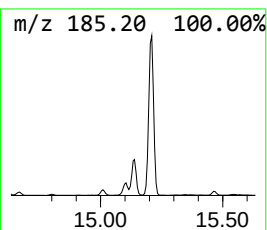
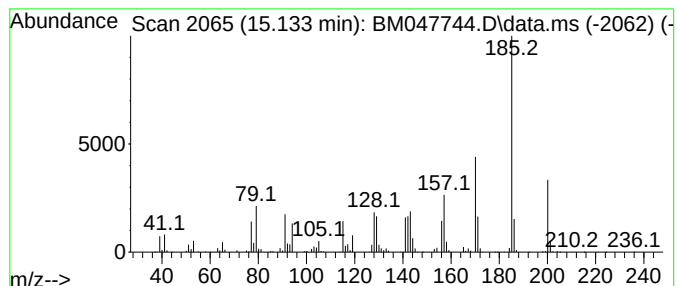
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BNA_M
ClientSampleId :
DDCA2

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Quant Title : SVOA CALIBRATION

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

Peak Number 51 9-Methyl-S-octahydroanthracene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.133	11.51 ng/ul	2247040	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Methyl-5-octahydroanthracene	200	C15H20	004948-51-0	70
2	1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	64
3	.alpha.-Corocalene	200	C15H20	020129-39-9	53
4	2-Acetyl-6-methoxynaphthalene	200	C13H12O2	003900-45-6	52
5	5-Hydroxy-4-nitroguaiacol	185	C7H7NO5	100960-04-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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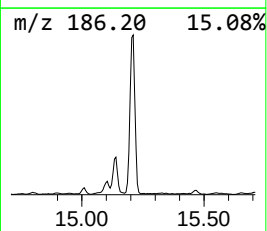
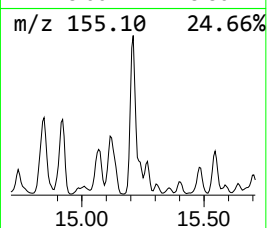
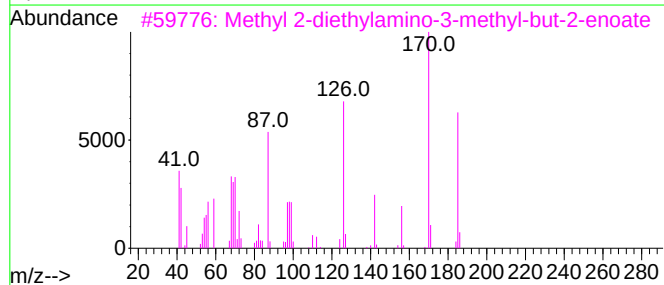
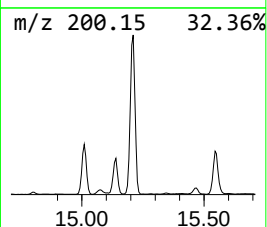
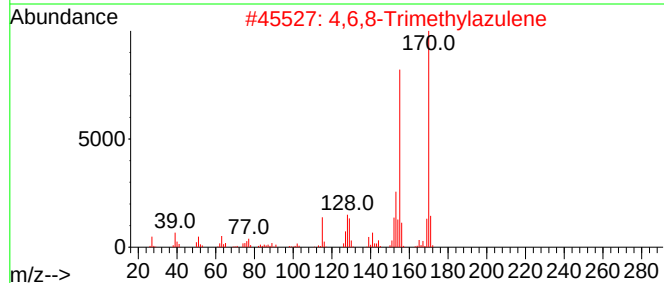
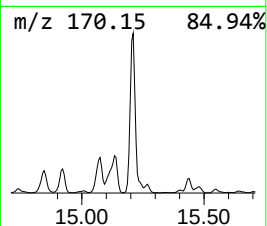
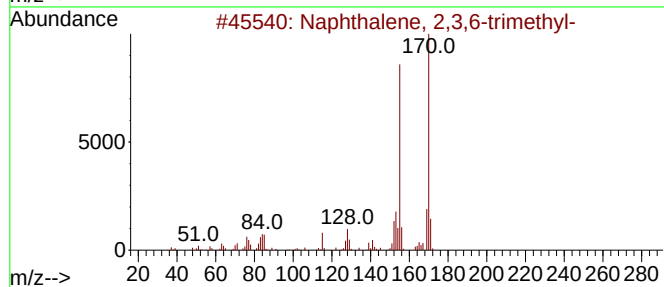
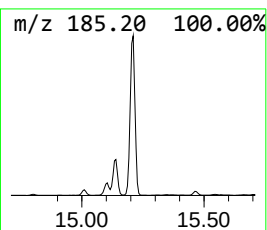
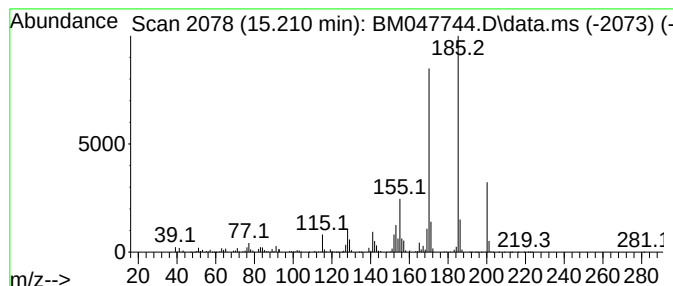
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 4,6,8-Trimethylazulene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	40.32 ng/ul	7871850	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	60
2			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
3			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	50
4			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	50
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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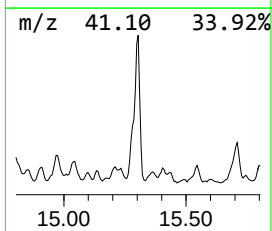
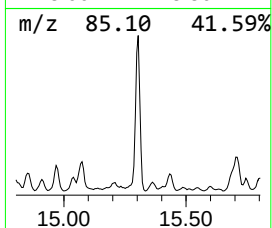
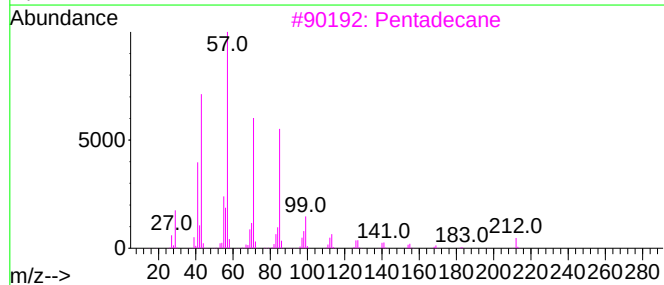
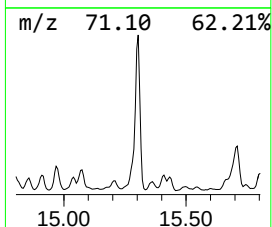
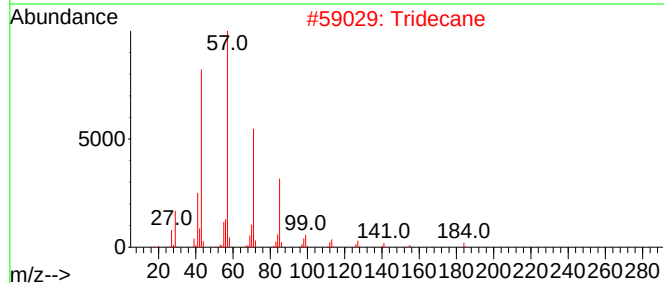
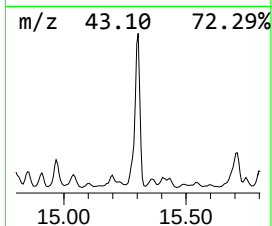
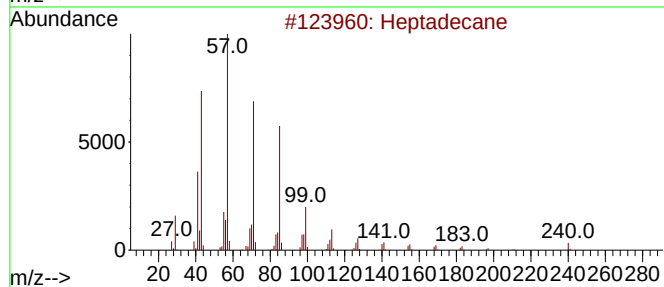
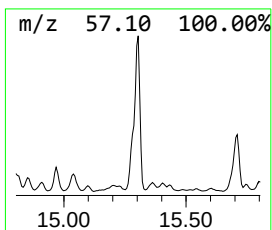
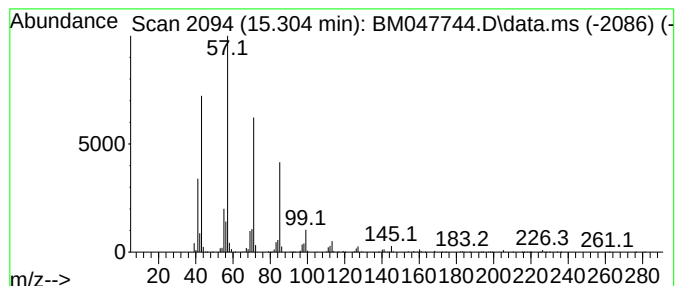
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.304	33.83 ng/ul	6605260	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	90
2			Tridecane	184	C13H28	000629-50-5	86
3			Pentadecane	212	C15H32	000629-62-9	86
4			Eicosane	282	C20H42	000112-95-8	86
5			Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

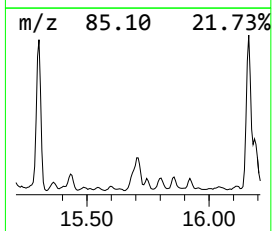
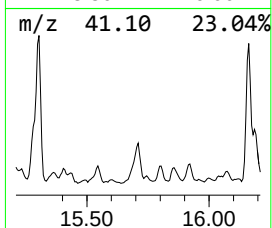
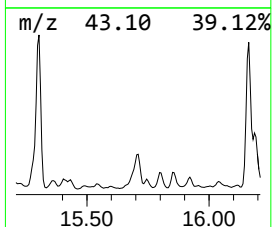
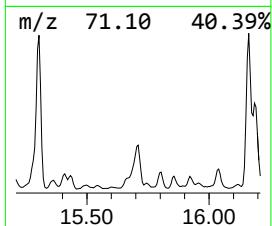
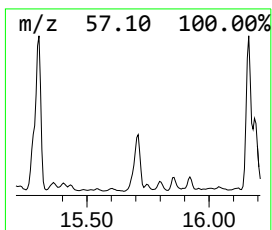
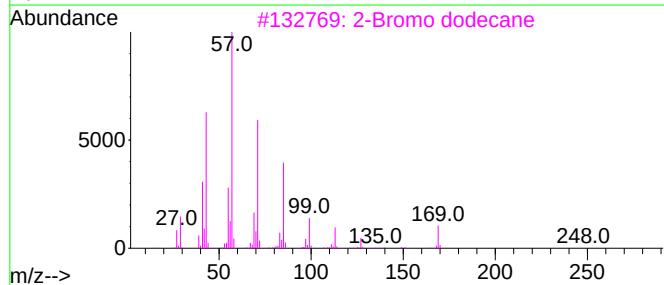
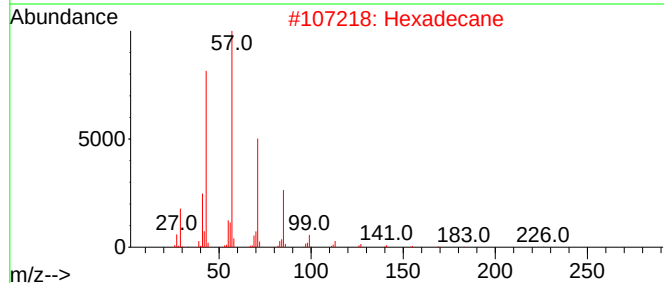
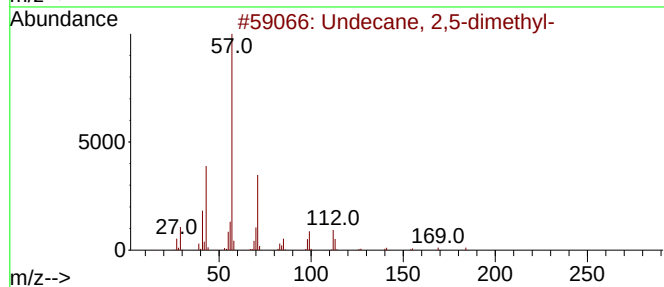
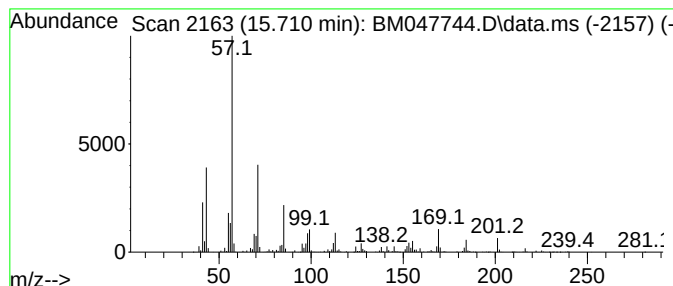
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.710	9.24 ng/ul	1803470	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	70
2	Hexadecane	226	C16H34	000544-76-3	70
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	68
4	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	58
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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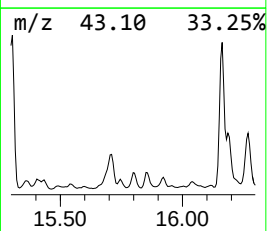
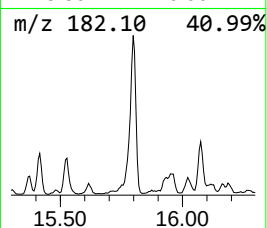
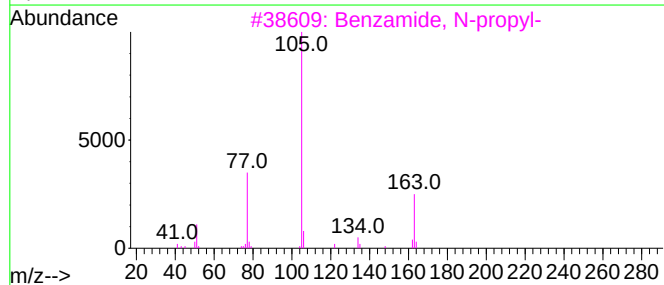
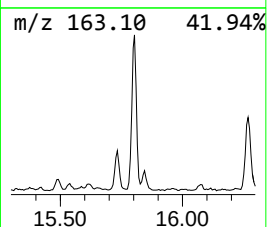
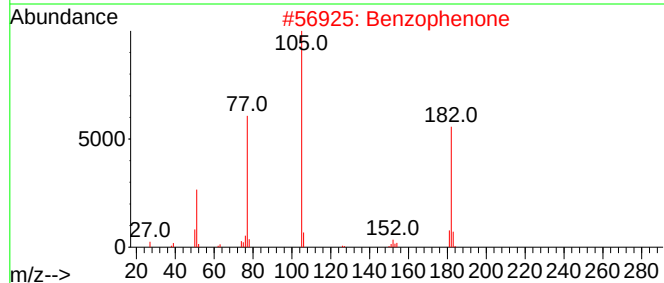
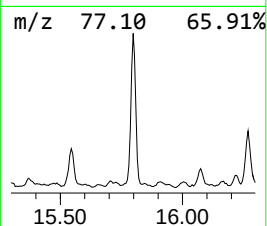
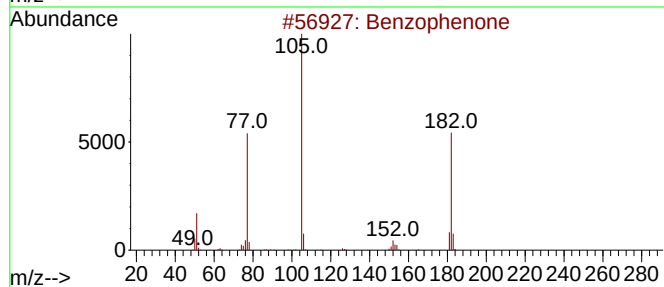
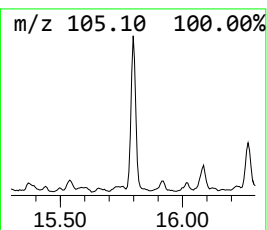
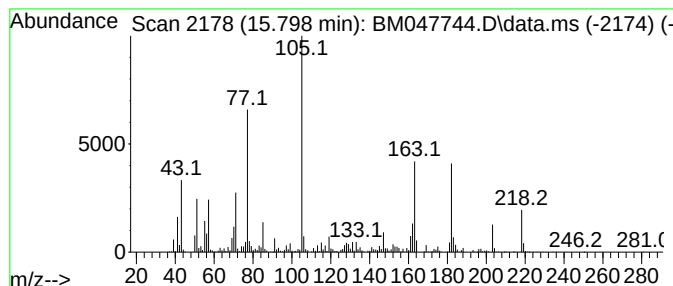
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Quant Title : SVOA CALIBRATION

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

Peak Number 55 Benzophenone Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.798	10.71 ng/ul	2091790	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	84
2			Benzophenone	182	C13H10O	000119-61-9	52
3			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	50
4			1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	41
5			Benzophenone	182	C13H10O	000119-61-9	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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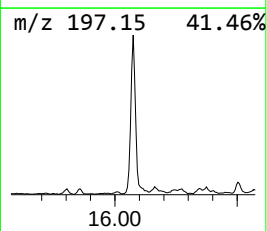
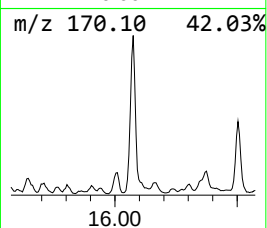
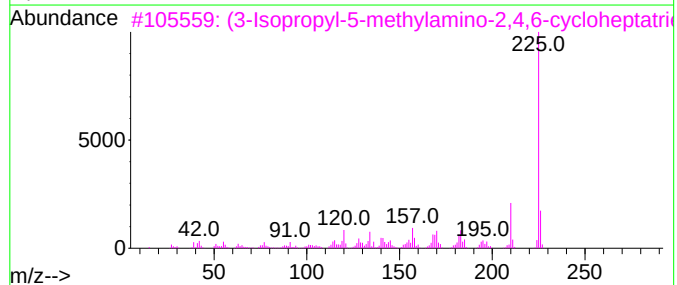
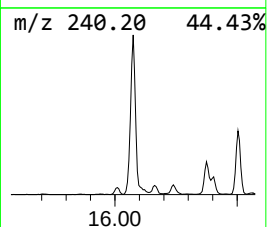
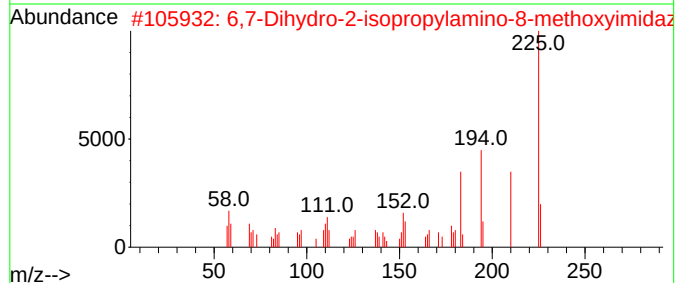
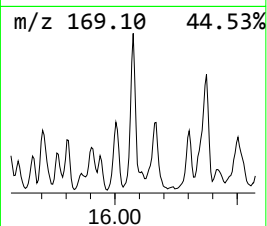
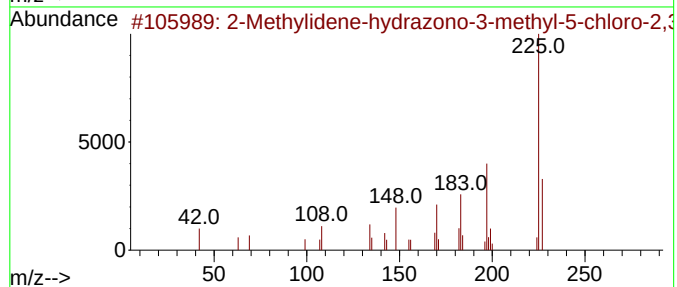
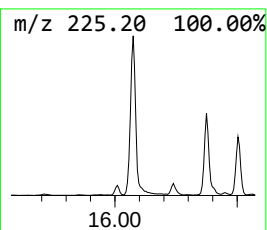
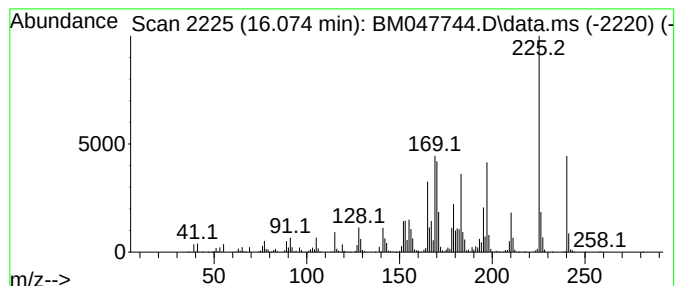
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 unknown-03 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.074	12.00 ng/ul	2716150	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	43		
2	6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	42		
3	(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	30		
4	benzoic acid, 2-amino-5-(butylth...	225	C11H15N02S	1000400-91-1	27		
5	Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
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Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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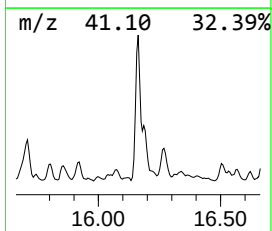
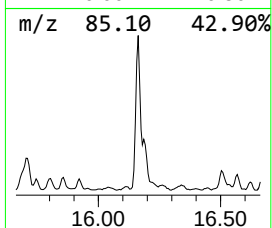
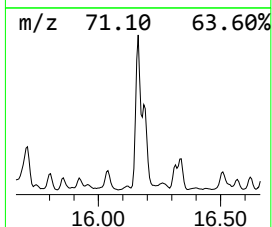
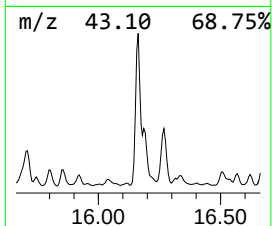
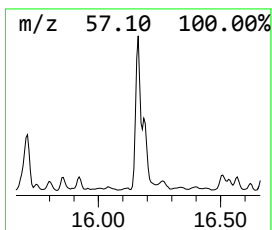
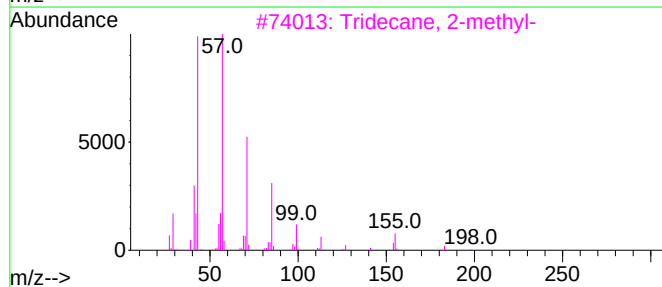
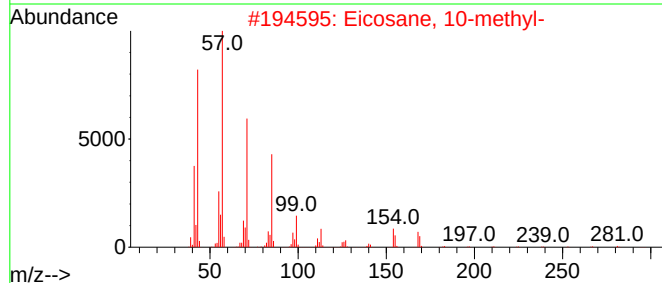
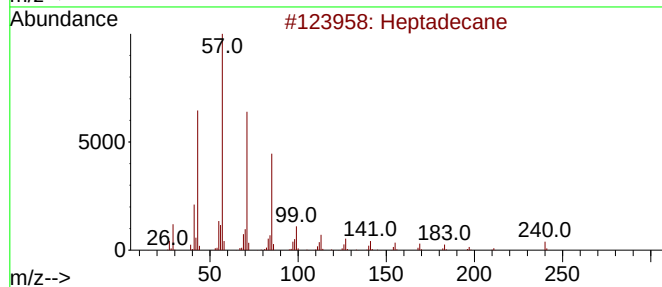
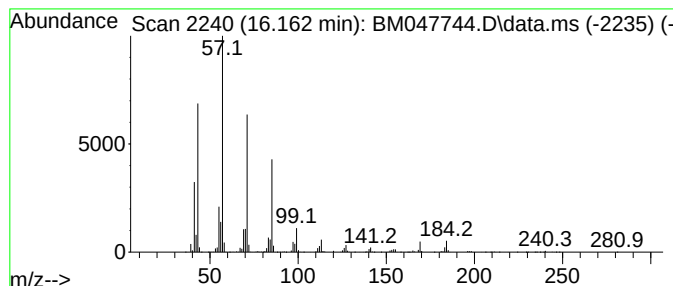
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.163	23.42 ng/ul	5299930	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	94
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
3			Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
4			Heneicosane	296	C21H44	000629-94-7	86
5			Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

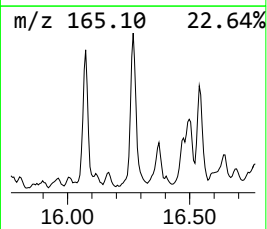
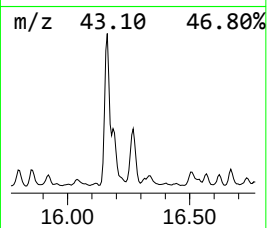
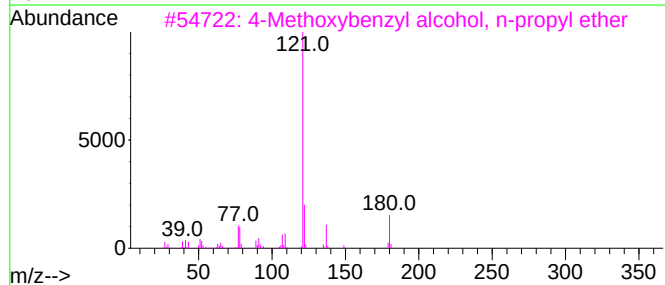
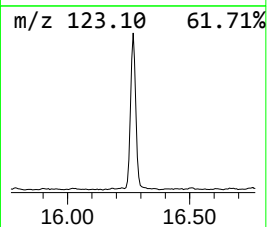
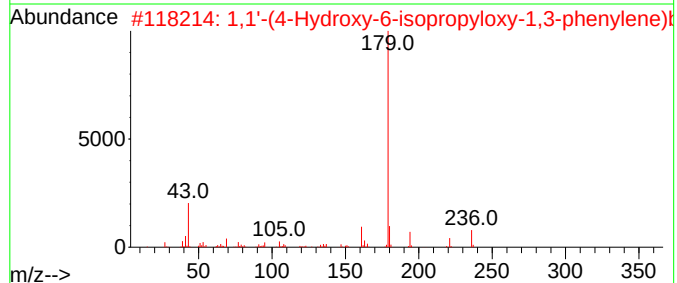
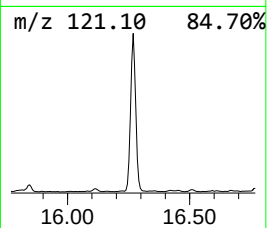
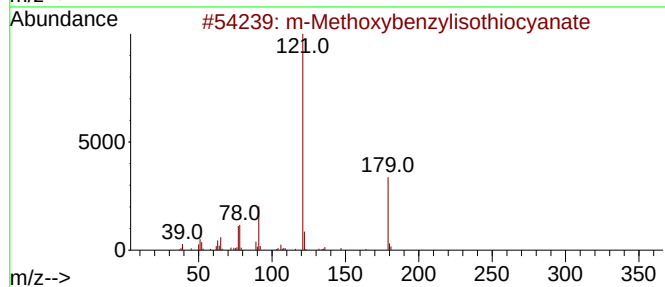
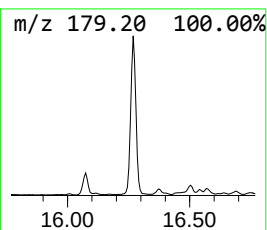
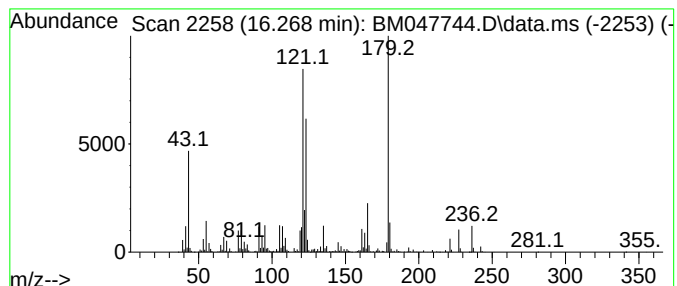
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 unknown-04 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.268	17.85 ng/ul	4038970	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Methoxybenzylisothiocyanate	179	C9H9NOS	075272-77-4	49
2			1,1'-(4-Hydroxy-6-isopropoxy-1...	236	C13H16O4	1000454-72-6	42
3			4-Methoxybenzyl alcohol, n-propy...	180	C11H16O2	1000394-78-1	27
4			4-Dimethylamino-2-methoxybenzald...	179	C10H13NO2	084562-48-1	25
5			Benzothiazole-6-carboxylic acid	179	C8H5NO2S	003622-35-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

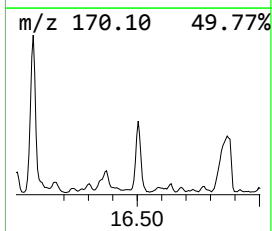
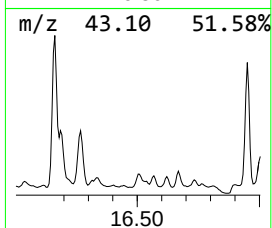
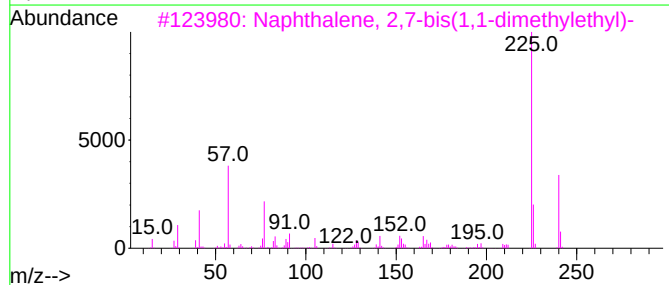
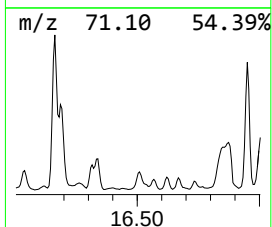
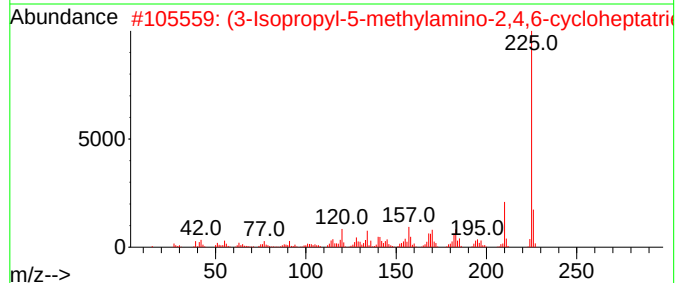
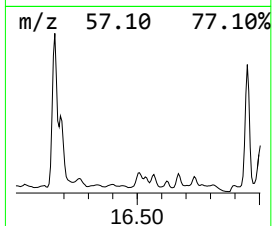
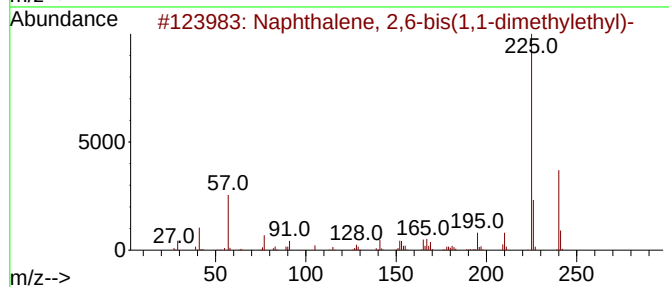
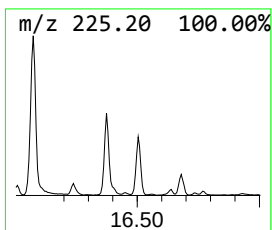
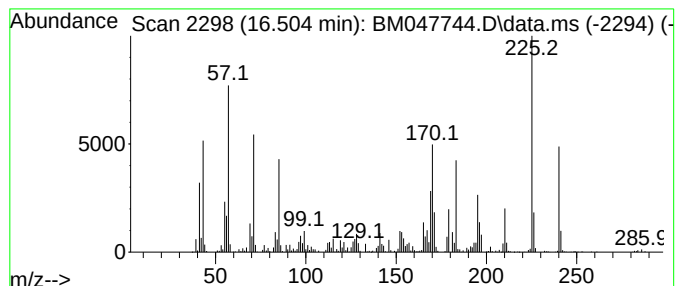
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 unknown-05 Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.504	6.26 ng/ul	1416910	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	27
2			(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	25
3			Naphthalene, 2,7-bis(1,1-dimethy...	240	C18H24	010275-58-8	22
4			1H-Pyrazole, 5-(4-ethylphenyl)-3...	240	C12H11F3N2	1000310-25-4	22
5			ONN-Azoxybenzene, 2,4,6-trimethyl-	240	C15H16N2O	016914-58-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

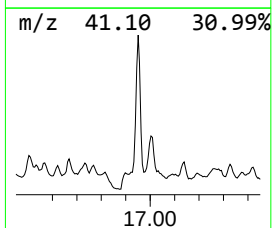
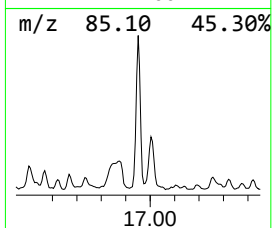
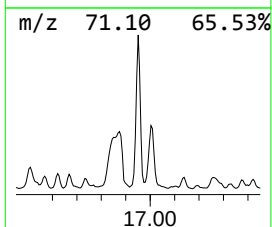
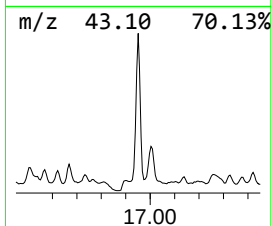
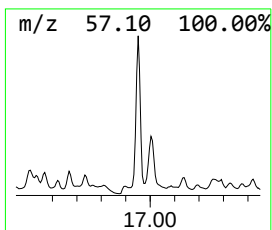
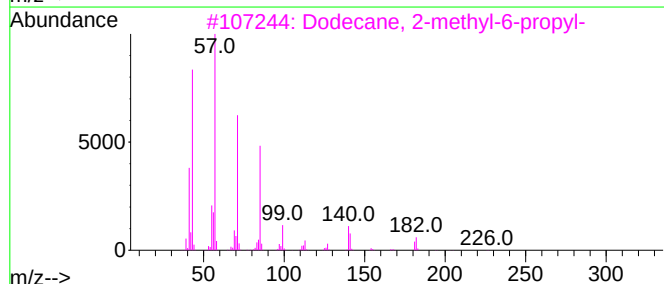
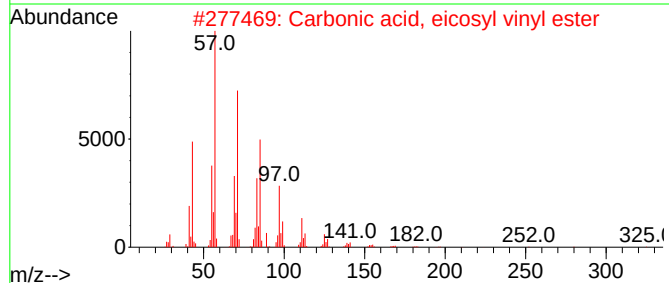
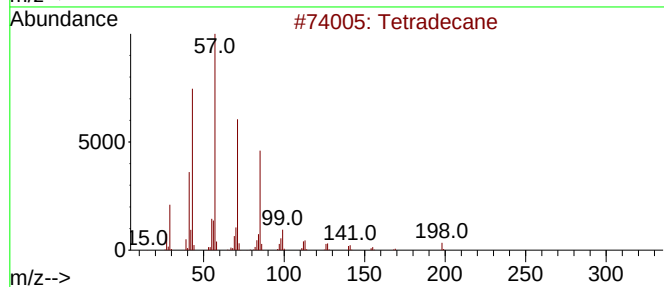
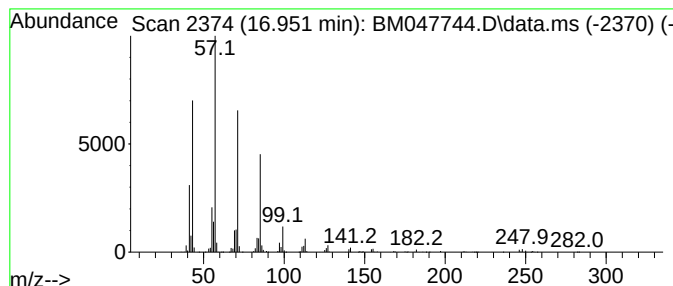
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 60 Carbonic acid, eicosyl vinyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.951	18.50 ng/ul	4186030	Phenanthrene-d10		17.186	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	86
2	Carbonic acid, eicosyl vinyl ester		368	C23H44O3	1000382-54-3	80
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	80
4	Nonane, 5-butyl-		184	C13H28	017312-63-9	64
5	Heptacosane		380	C27H56	000593-49-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

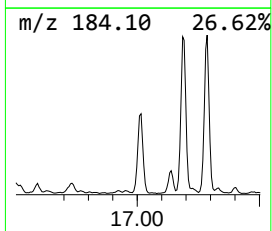
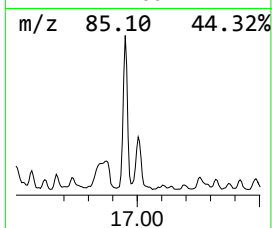
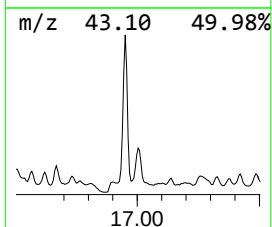
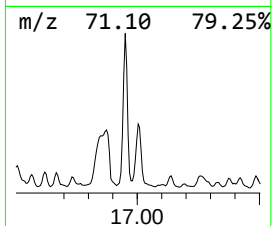
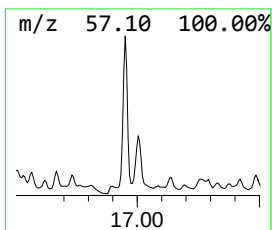
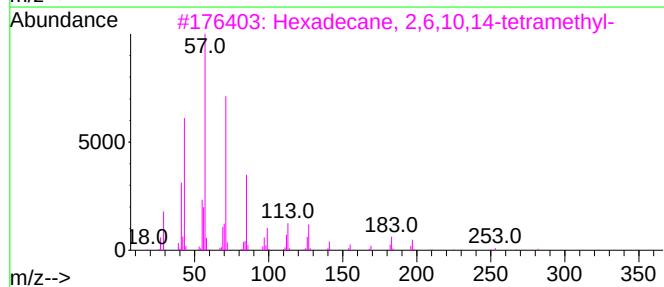
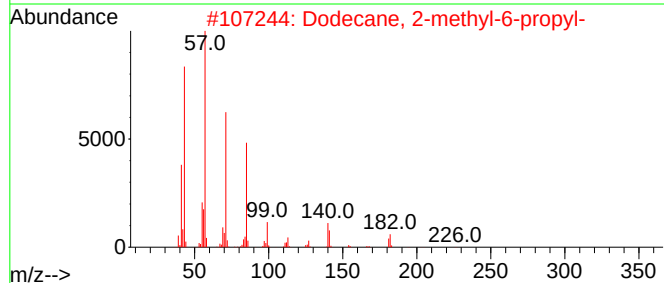
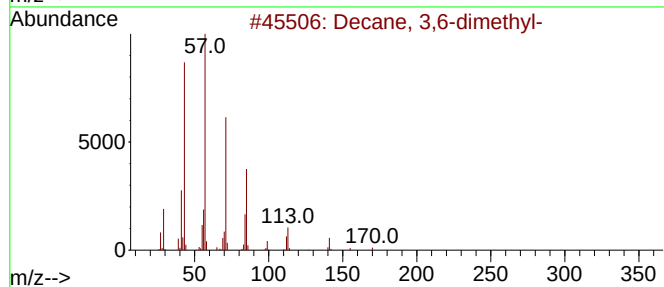
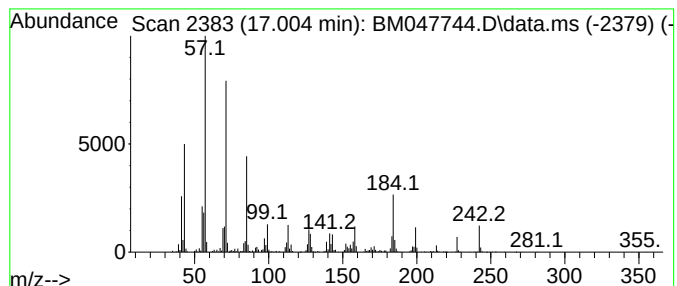
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.004	11.64 ng/ul	2634280	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	70
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	70
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	68
4	Eicosane	282	C20H42	000112-95-8	64
5	Tetracosane	338	C24H50	000646-31-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

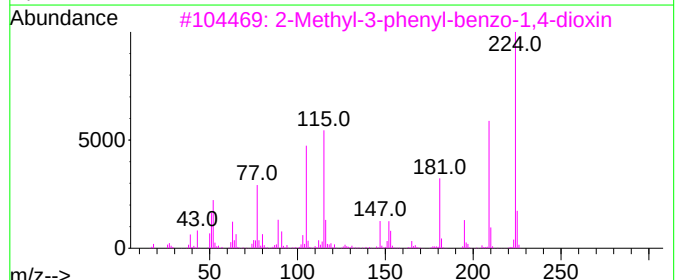
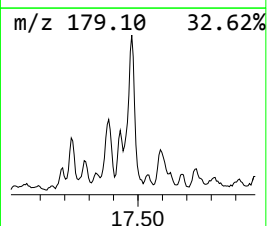
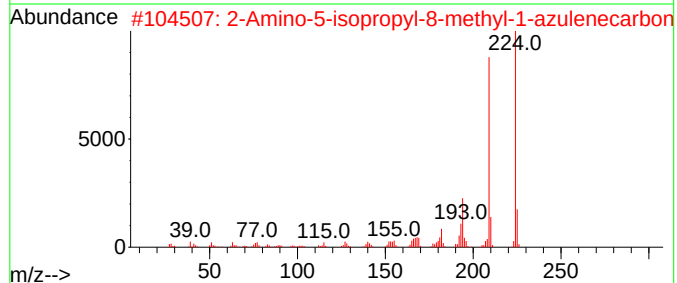
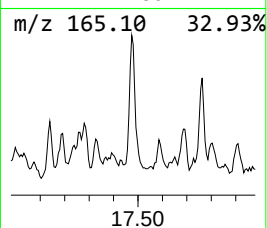
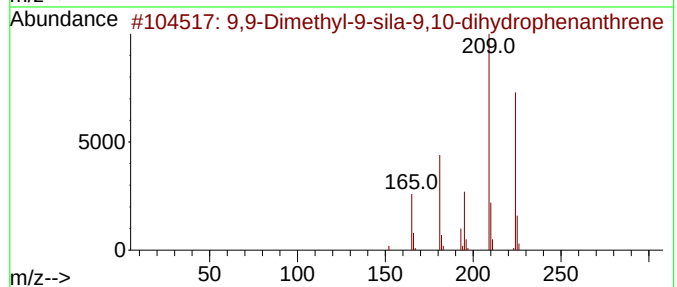
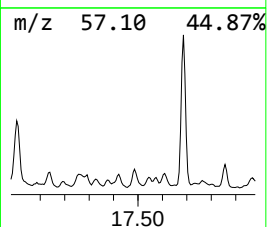
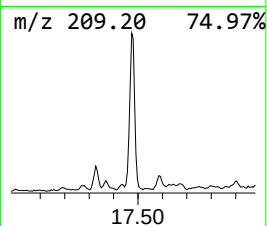
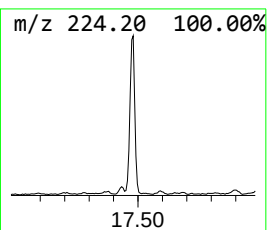
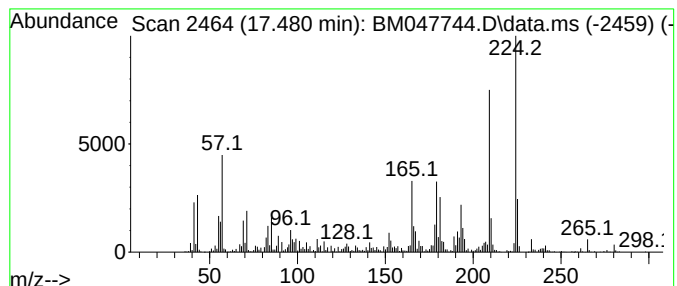
Instrument :
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ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 62 9,9-Dimethyl-9-sila-9,10-di... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.480	8.69 ng/ul	1966010	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si	187328-55-8	81
2	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	62
3	2-Methyl-3-phenyl-benzo-1,4-dioxin	224	C15H12O2	079792-91-9	58
4	7-Methyl-6,8-bis(methylthio)pyrr...	224	C10H12N2S2	084201-40-1	52
5	1,2,3,4-Tetrahydronaphthalen-1,2...	224	C12H16O4	1000125-87-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

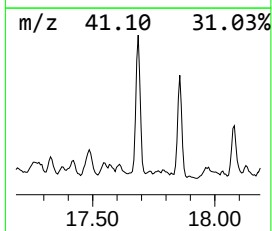
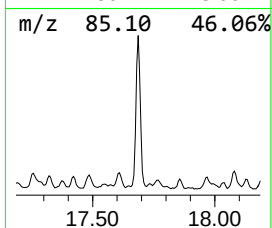
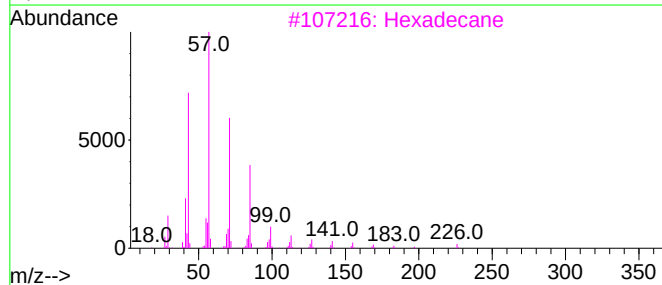
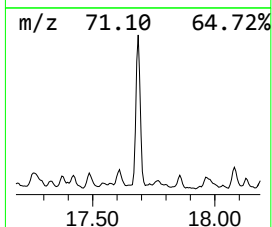
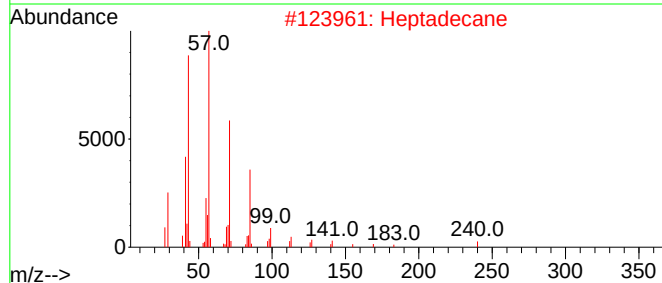
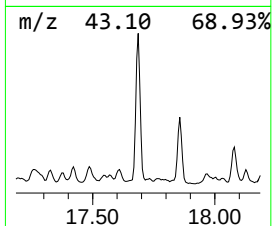
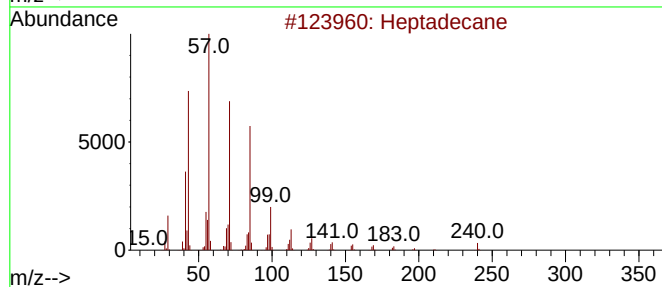
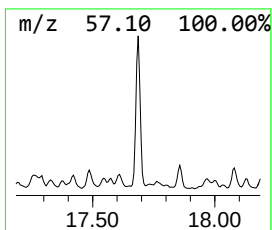
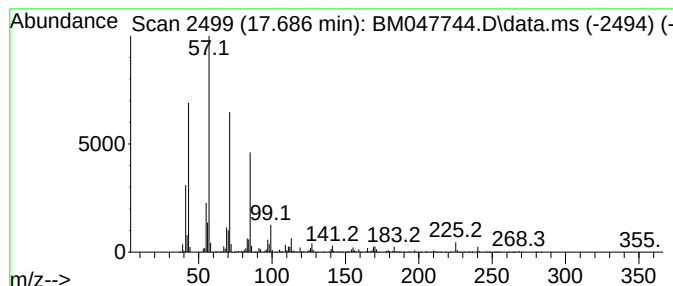
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.686	14.48 ng/ul	3277300	Phenanthrene-d10		17.186	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Heptadecane		240	C17H36	000629-78-7	94
3	Hexadecane		226	C16H34	000544-76-3	93
4	Nonadecane		268	C19H40	000629-92-5	91
5	3-Ethyl-2,6,10-trimethylundecane		226	C16H34	1000432-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

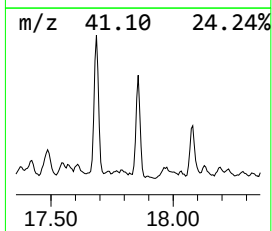
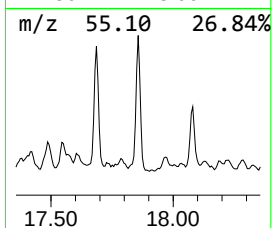
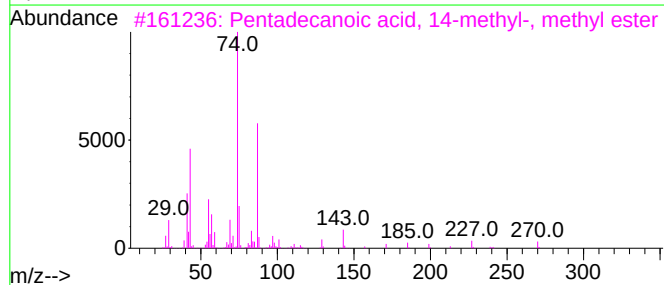
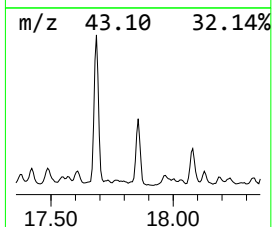
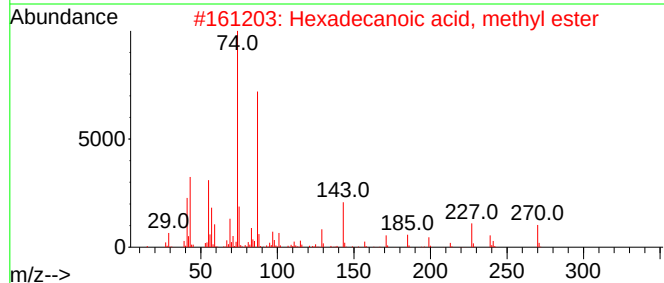
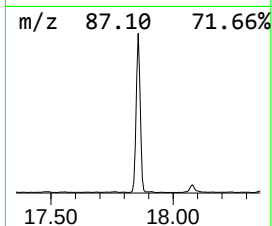
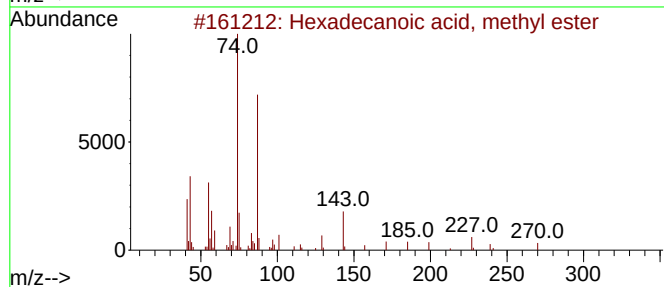
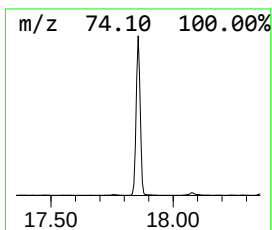
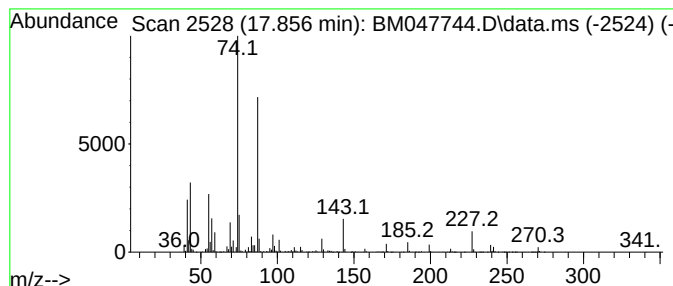
Instrument :
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ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 64 Hexadecanoic acid, methyl e... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.857	11.29 ng/ul	2555280	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
3	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

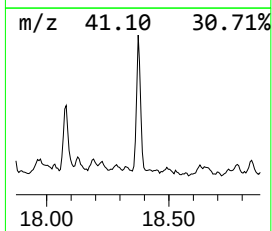
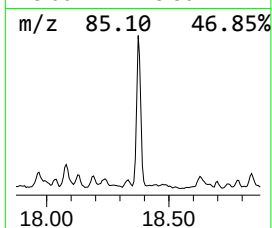
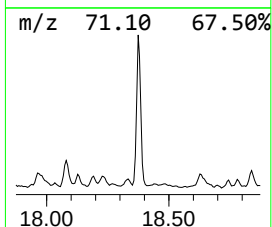
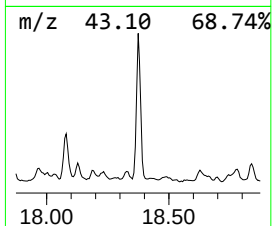
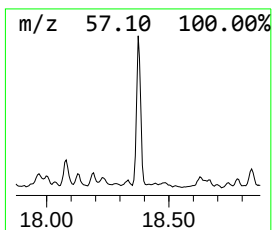
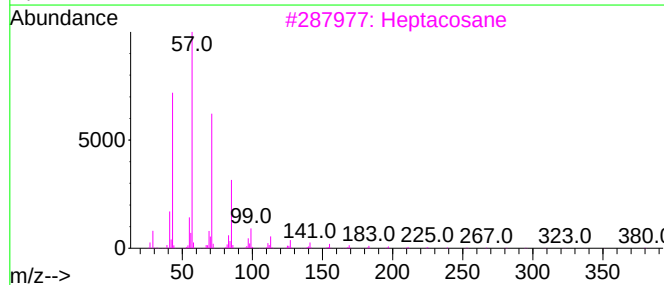
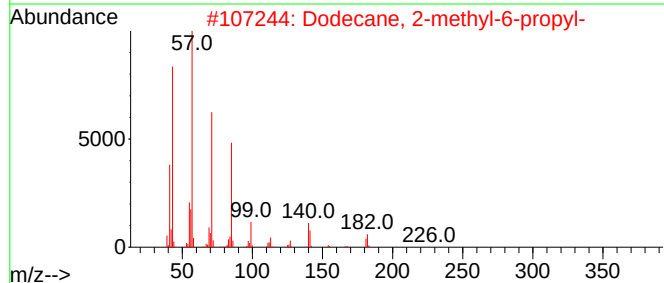
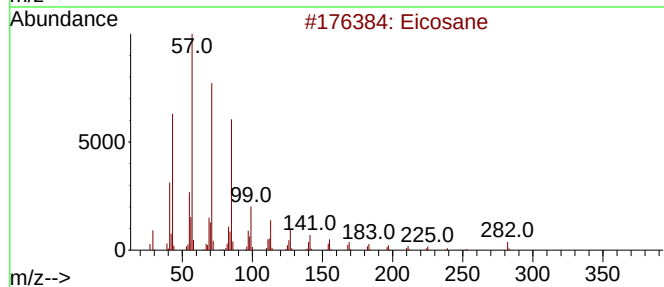
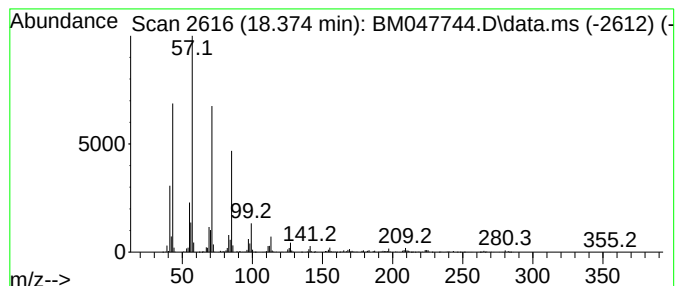
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Quant Title : SVOA CALIBRATION

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

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.374	11.39 ng/ul	2577950	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	97
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
3		Heptacosane	380	C27H56	000593-49-7	91
4		Nonadecane	268	C19H40	000629-92-5	91
5		Nonadecane	268	C19H40	000629-92-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

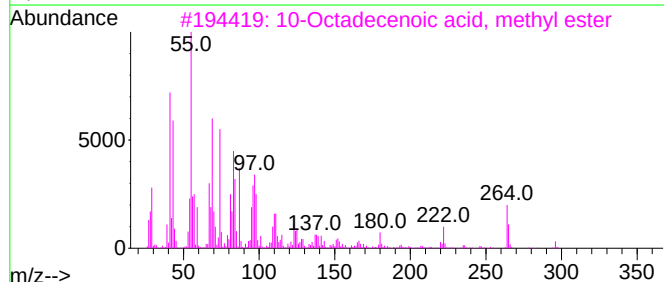
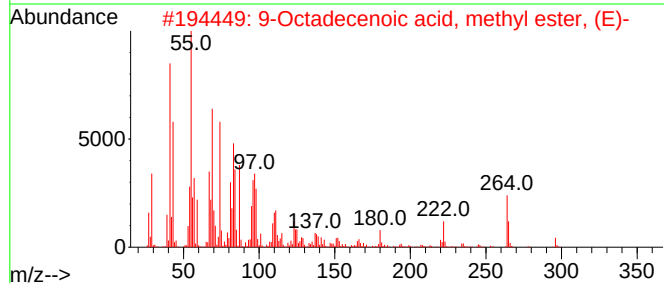
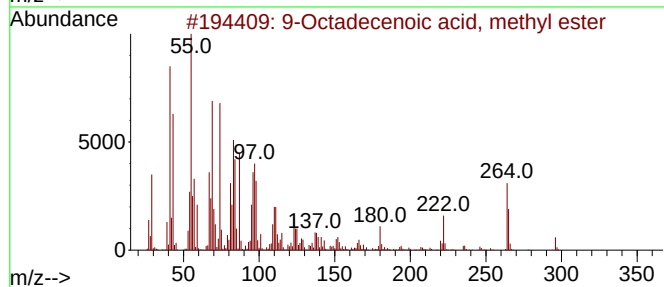
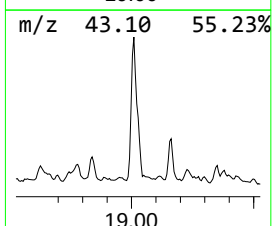
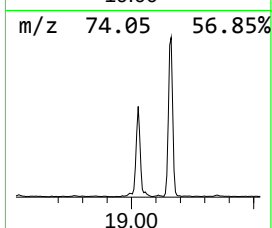
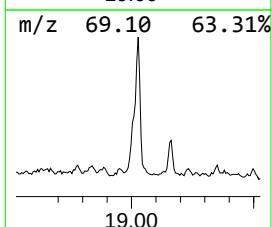
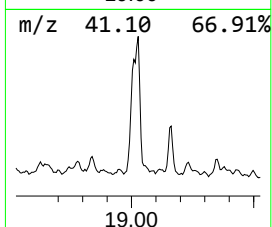
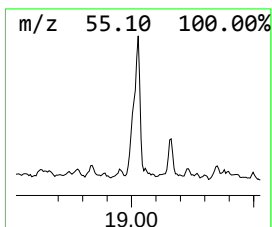
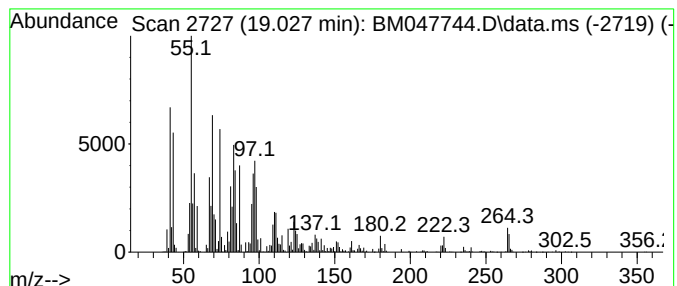
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 67 9-Octadecenoic acid, methyl... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.027	17.70 ng/ul	4005110	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
2	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
3	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4	6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	99
5	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

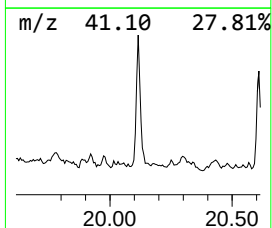
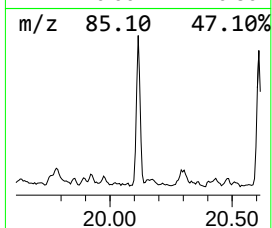
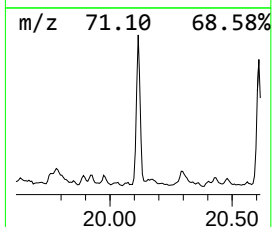
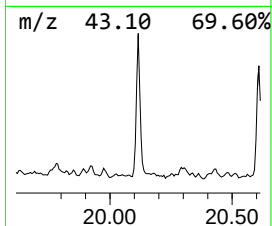
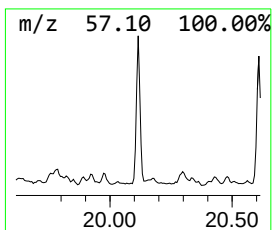
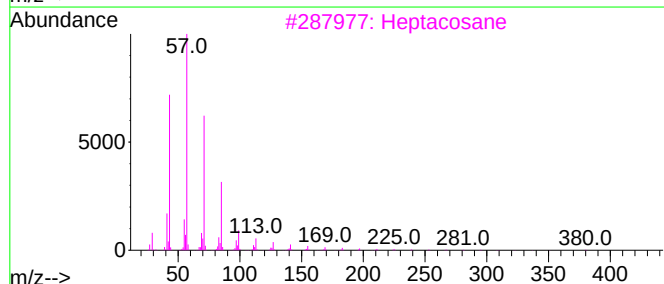
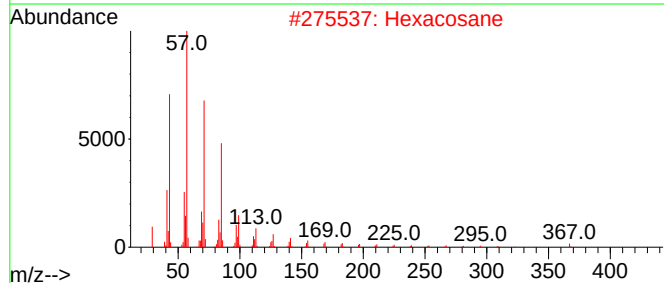
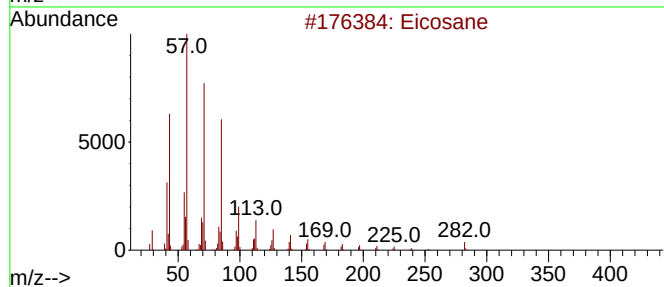
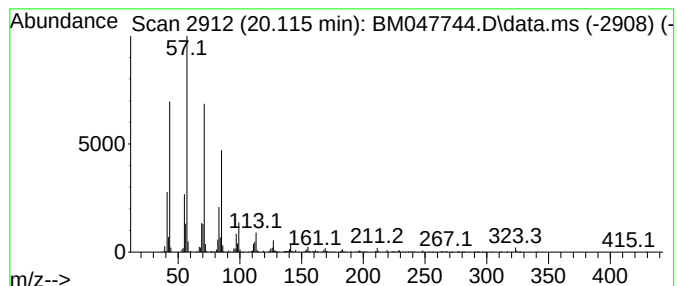
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.115	8.89 ng/ul	1782650	Chrysene-d12	21.403	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Hexacosane	366	C26H54	000630-01-3	87
3	Heptacosane	380	C27H56	000593-49-7	87
4	Octacosane	394	C28H58	000630-02-4	87
5	Pentadecane	212	C15H32	000629-62-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCA2

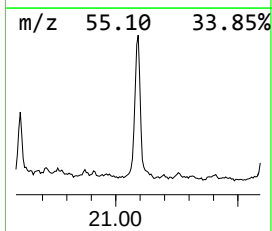
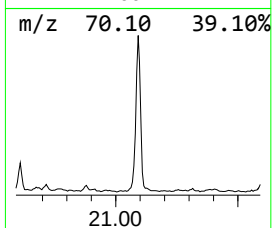
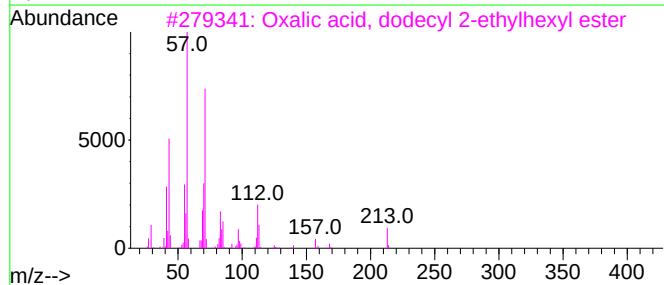
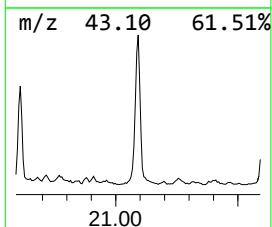
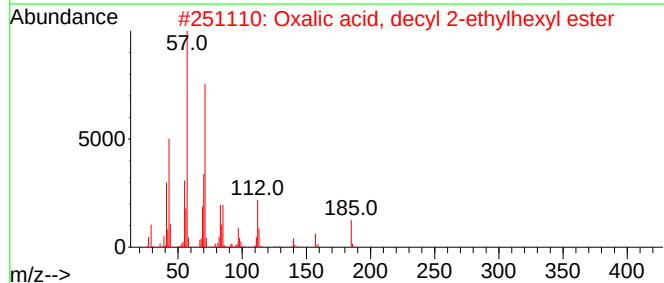
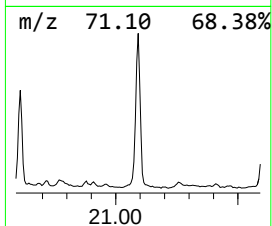
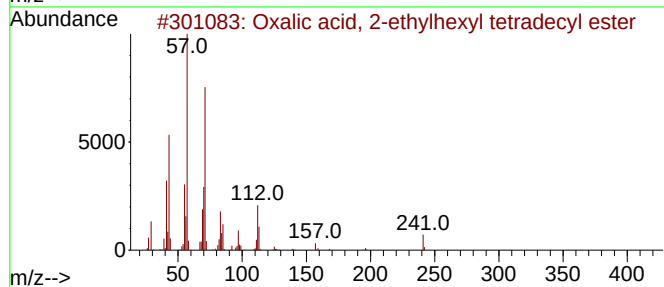
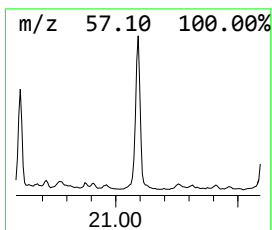
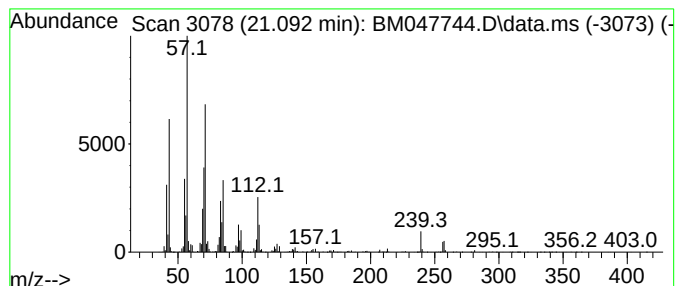
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Quant Title : SVOA CALIBRATION

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

Peak Number 69 Oxalic acid, 2-ethylhexyl t... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	13.56 ng/ul	2718210	Chrysene-d12	21.403

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	83
2			Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	80
3			Oxalic acid, dodecyl 2-ethylhexy...	370	C22H42O4	1000309-39-5	74
4			Carbonic acid, 2-ethylhexyl trid...	356	C22H44O3	1000383-14-0	72
5			Oxalic acid, 2-ethylhexyl isohex...	286	C16H30O4	1000309-38-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047744.D
Acq On : 01 Oct 2024 15:43
Operator : RC/JU
Sample : P4018-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

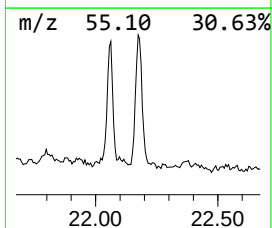
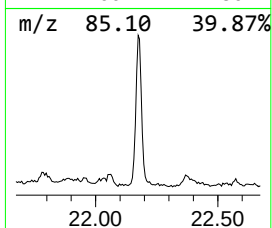
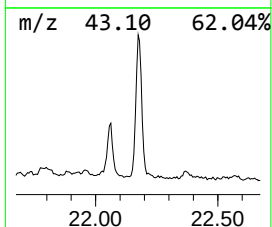
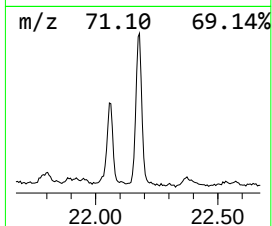
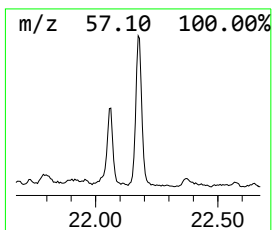
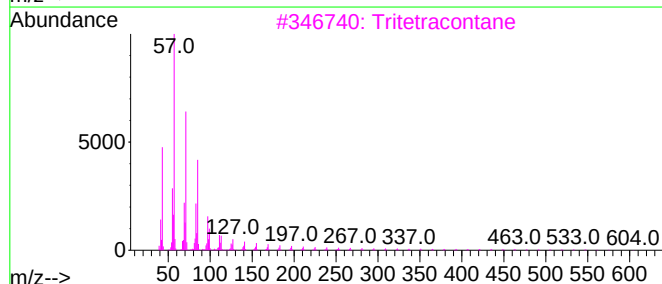
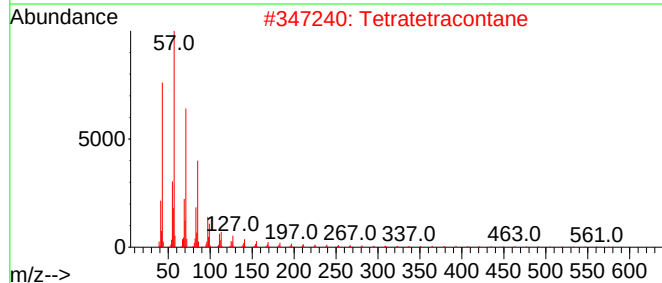
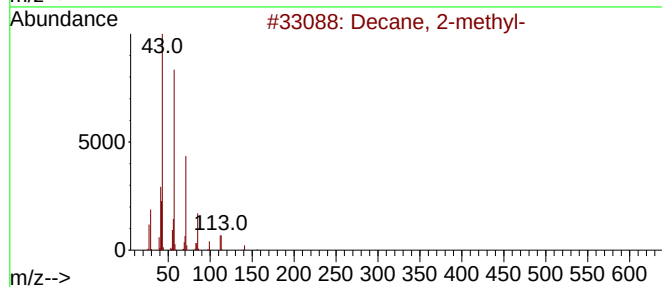
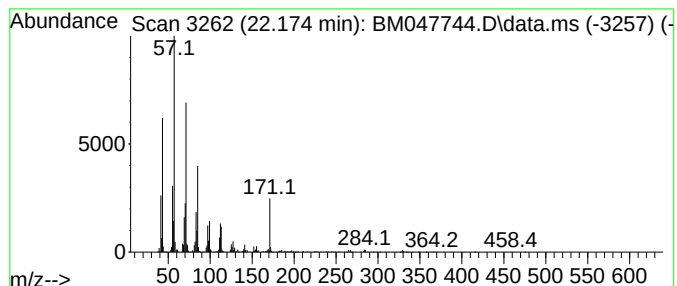
Instrument :
BNA_M
ClientSampleId :
DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.174	6.43 ng/ul	1288780	Chrysene-d12	21.403	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2-methyl-	156	C11H24	006975-98-0	76
2	Tetratetracontane	619	C44H90	007098-22-8	68
3	Tritetracontane	605	C43H88	007098-21-7	64
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	64
5	Decane, 5-propyl-	184	C13H28	017312-62-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
 Data File : BM047744.D
 Acq On : 01 Oct 2024 15:43
 Operator : RC/JU
 Sample : P4018-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDCA2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.834	22.2	ng/ul	4992190	1	7.810	4492900	20.0
Hexylene glycol	6.134	8.0	ng/ul	1807100	1	7.810	4492900	20.0
Hexyl octyl ether	6.316	8.3	ng/ul	1864360	1	7.810	4492900	20.0
(DEL) Alkane: S...	6.381	6.4	ng/ul	1442850	1	7.810	4492900	20.0
(DEL) Alkane: C...	6.457	7.9	ng/ul	1764970	1	7.810	4492900	20.0
(DEL) Alkane: S...	6.816	10.6	ng/ul	2381080	1	7.810	4492900	20.0
(DEL) Alkane: S...	6.869	11.1	ng/ul	2498240	1	7.810	4492900	20.0
Benzene, 1-ethy...	6.904	9.3	ng/ul	2093690	1	7.810	4492900	20.0
Mesitylene	7.040	9.5	ng/ul	2128700	1	7.810	4492900	20.0
Benzene, 1-ethy...	7.204	8.6	ng/ul	1921730	1	7.810	4492900	20.0
(DEL) Alkane: S...	7.445	74.2	ng/ul	16660200	1	7.810	4492900	20.0
Oxalic acid, is...	7.739	9.0	ng/ul	2013420	1	7.810	4492900	20.0
1-Hexanol, 2-et...	7.887	63.2	ng/ul	14190500	1	7.810	4492900	20.0
1,1-Hexylenedio...	8.039	23.8	ng/ul	5353290	1	7.810	4492900	20.0
(DEL) Alkane: C...	8.110	6.3	ng/ul	1404640	1	7.810	4492900	20.0
(DEL) Alkane: S...	8.351	15.9	ng/ul	3576280	1	7.810	4492900	20.0
(DEL) Alkane: S...	8.410	7.5	ng/ul	1692870	1	7.810	4492900	20.0
Benzene, 1,2-di...	8.457	9.9	ng/ul	2214070	1	7.810	4492900	20.0
(DEL) Alkane: S...	8.586	12.1	ng/ul	2708110	1	7.810	4492900	20.0
Benzene, 1-meth...	8.616	6.6	ng/ul	1489990	1	7.810	4492900	20.0
(DEL) Alkane: S...	9.051	50.7	ng/ul	11393400	1	7.810	4492900	20.0
Hexanoic acid, ...	9.175	27.3	ng/ul	6128770	1	7.810	4492900	20.0
3-Heptanol, 4-m...	10.492	9.1	ng/ul	4089110	2	10.604	8980330	20.0
(DEL) Alkane: S...	10.986	7.2	ng/ul	3242090	2	10.604	8980330	20.0
Benzene, 1,3,5-...	11.704	14.2	ng/ul	6395080	2	10.604	8980330	20.0
(DEL) Alkane: S...	12.033	9.2	ng/ul	4137400	2	10.604	8980330	20.0
3-Trifluoroacet...	12.680	20.3	ng/ul	3953870	3	14.445	3905050	20.0
(DEL) Alkane: S...	13.280	22.3	ng/ul	4357120	3	14.445	3905050	20.0
unknown-01	13.716	6.9	ng/ul	1350620	3	14.445	3905050	20.0
Naphthalene, 2,...	13.769	8.6	ng/ul	1687830	3	14.445	3905050	20.0
unknown-02	14.080	10.1	ng/ul	1978240	3	14.445	3905050	20.0
(DEL) Alkane: S...	14.363	132.8	ng/ul	25924800	3	14.445	3905050	20.0
(DEL) Alkane: S...	14.763	17.5	ng/ul	3409310	3	14.445	3905050	20.0
Naphthalene, 2,...	14.916	8.1	ng/ul	1574440	3	14.445	3905050	20.0
(DEL) Alkane: S...	14.974	7.2	ng/ul	1405180	3	14.445	3905050	20.0
4b,5,6,7,8,8a,9...	15.010	6.8	ng/ul	1328330	3	14.445	3905050	20.0
9-Methyl-S-octa...	15.133	11.5	ng/ul	2247040	3	14.445	3905050	20.0
4,6,8-Trimethyl...	15.210	40.3	ng/ul	7871850	3	14.445	3905050	20.0
(DEL) Alkane: S...	15.304	33.8	ng/ul	6605260	3	14.445	3905050	20.0
(DEL) Alkane: S...	15.710	9.2	ng/ul	1803470	3	14.445	3905050	20.0
Benzophenone	15.798	10.7	ng/ul	2091790	3	14.445	3905050	20.0
unknown-03	16.074	12.0	ng/ul	2716150	4	17.186	4526470	20.0
(DEL) Alkane: S...	16.163	23.4	ng/ul	5299930	4	17.186	4526470	20.0
unknown-04	16.268	17.9	ng/ul	4038970	4	17.186	4526470	20.0
unknown-05	16.504	6.3	ng/ul	1416910	4	17.186	4526470	20.0
Carbonic acid, ...	16.951	18.5	ng/ul	4186030	4	17.186	4526470	20.0
(DEL) Alkane: S...	17.004	11.6	ng/ul	2634280	4	17.186	4526470	20.0
9,9-Dimethyl-9-...	17.480	8.7	ng/ul	1966010	4	17.186	4526470	20.0
(DEL) Alkane: S...	17.686	14.5	ng/ul	3277300	4	17.186	4526470	20.0
Hexadecanoic ac...	17.857	11.3	ng/ul	2555280	4	17.186	4526470	20.0
(DEL) Alkane: S...	18.374	11.4	ng/ul	2577950	4	17.186	4526470	20.0
9-Octadecenoic ...	19.027	17.7	ng/ul	4005110	4	17.186	4526470	20.0

(DEL) Alkane: S...	20.115	8.9	ng/ul	1782650	5	21.403	4009660	20.0
Oxalic acid, 2-...	21.092	13.6	ng/ul	2718210	5	21.403	4009660	20.0
(DEL) Alkane: S...	22.174	6.4	ng/ul	1288780	5	21.403	4009660	20.0

Instrument :
BNA_M
ClientSampleId :
DDCA2