

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
 Data File : BM045843.D
 Acq On : 05 Jun 2024 09:00
 Operator : MA/JU
 Sample : P2586-03
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4D16

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-BM053124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM045843.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.052	8	11	20	rVB	78310	113649	2.51%	0.158%
2	3.463	78	81	103	rBV	764781	1557572	34.35%	2.161%
3	3.746	126	129	135	rVB2	27296	38070	0.84%	0.053%
4	4.169	197	201	208	rVB3	23246	43920	0.97%	0.061%
5	4.252	213	215	220	rVB4	13106	16448	0.36%	0.023%
6	4.987	335	340	344	rBV7	6279	11116	0.25%	0.015%
7	5.651	448	453	460	rBV3	25734	55997	1.23%	0.078%
8	5.757	469	471	478	rVB8	10332	18630	0.41%	0.026%
9	6.722	628	635	644	rBV	1433274	2393090	52.78%	3.321%
10	6.828	647	653	665	rVV	1040169	1660012	36.61%	2.303%
11	7.028	680	687	696	rBV	1469690	2298774	50.70%	3.190%
12	7.463	754	761	769	rBV	1352122	2042512	45.05%	2.834%
13	7.575	776	780	785	rBV4	14817	27729	0.61%	0.038%
14	7.616	785	787	792	rVV6	10121	15645	0.35%	0.022%
15	8.234	884	892	899	rBV	1727201	2770143	61.09%	3.844%
16	8.292	899	902	911	rVB2	71893	124515	2.75%	0.173%
17	8.628	951	959	971	rBV	1461808	2313607	51.03%	3.210%
18	8.804	985	989	994	rBV5	22168	31899	0.70%	0.044%
19	9.239	1053	1063	1069	rVV	119173	212897	4.70%	0.295%
20	9.339	1073	1080	1099	rVB	1230929	2010740	44.35%	2.790%
21	9.687	1134	1139	1143	rBV6	13021	25937	0.57%	0.036%
22	9.886	1165	1173	1187	rBV	1628657	2803646	61.83%	3.890%
23	10.234	1222	1232	1247	rBV	1716545	2813844	62.06%	3.904%
24	10.410	1254	1262	1280	rBV	1245946	2424635	53.47%	3.364%
25	10.857	1333	1338	1347	rBV3	35030	68622	1.51%	0.095%
26	11.028	1358	1367	1379	rBV	134617	308487	6.80%	0.428%
27	11.122	1379	1383	1391	rVB9	10402	26238	0.58%	0.036%
28	11.681	1476	1478	1487	rVB10	9077	19517	0.43%	0.027%
29	11.833	1497	1504	1510	rVB	32981	57086	1.26%	0.079%
30	11.998	1526	1532	1535	rBV4	12813	23296	0.51%	0.032%
31	12.133	1552	1555	1560	rVB6	5816	9678	0.21%	0.013%
32	12.869	1676	1680	1684	rBV3	25335	35724	0.79%	0.050%
33	12.945	1689	1693	1697	rBV4	9494	14435	0.32%	0.020%
34	12.992	1697	1701	1706	rVB4	17012	25324	0.56%	0.035%
35	13.080	1711	1716	1721	rVB8	10708	24110	0.53%	0.033%
36	13.327	1750	1758	1768	rVV	69457	174401	3.85%	0.242%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
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37	13.427	1768	1775	1778	rVV6	47629	85040	1.88%	0.118%
38	13.463	1778	1781	1788	rVB3	42800	60804	1.34%	0.084%
39	13.557	1788	1797	1804	rBV	2539605	3449019	76.07%	4.786%
40	13.616	1805	1807	1810	rVV4	13979	16506	0.36%	0.023%

41	13.798	1830	1838	1848	rBV2	2752481	4137858	91.26%	5.742%
42	13.880	1848	1852	1855	rVV6	30442	43715	0.96%	0.061%
43	13.957	1859	1865	1871	rVB4	62011	94162	2.08%	0.131%
44	14.027	1871	1877	1883	rVB9	16439	34635	0.76%	0.048%
45	14.104	1883	1890	1897	rBV	2244527	3163038	69.76%	4.389%

46	14.186	1899	1904	1907	rVB7	13288	18658	0.41%	0.026%
47	14.404	1927	1941	1959	rBV	899403	1654723	36.49%	2.296%
48	14.539	1959	1964	1968	rVV8	15171	32425	0.72%	0.045%
49	14.592	1970	1973	1980	rVB8	13616	23715	0.52%	0.033%
50	14.698	1985	1991	1998	rVV6	14967	30105	0.66%	0.042%

51	14.792	2003	2007	2009	rBV4	11513	13439	0.30%	0.019%
52	14.851	2014	2017	2021	rVB6	7708	11930	0.26%	0.017%
53	14.921	2024	2029	2033	rBV6	12354	18981	0.42%	0.026%
54	14.980	2033	2039	2042	rBV7	13156	19565	0.43%	0.027%
55	15.045	2044	2050	2054	rBV8	14490	26299	0.58%	0.036%

56	15.104	2054	2060	2073	rVB	3257718	4534190	100.00%	6.291%
57	15.245	2073	2084	2091	rBV	728827	1041821	22.98%	1.446%
58	15.357	2097	2103	2104	rBV6	15199	26383	0.58%	0.037%
59	15.445	2114	2118	2120	rBV4	11482	16689	0.37%	0.023%
60	15.498	2124	2127	2131	rVB5	16642	22135	0.49%	0.031%

61	15.557	2131	2137	2140	rBV7	38287	65867	1.45%	0.091%
62	15.598	2141	2144	2153	rVV6	33651	91217	2.01%	0.127%
63	15.668	2153	2156	2157	rVV3	10919	11149	0.25%	0.015%
64	15.704	2158	2162	2165	rVV3	36409	51347	1.13%	0.071%
65	15.757	2169	2171	2175	rVB5	10425	10943	0.24%	0.015%

66	15.845	2181	2186	2190	rBV3	49433	72811	1.61%	0.101%
67	15.886	2192	2193	2200	rVB7	11625	11797	0.26%	0.016%
68	16.033	2216	2218	2223	rVB6	10315	11959	0.26%	0.017%
69	16.263	2252	2257	2260	rBV7	10801	15830	0.35%	0.022%
70	16.316	2261	2266	2271	rBV7	25779	41648	0.92%	0.058%

71	16.386	2275	2278	2281	rBV4	18004	23413	0.52%	0.032%
72	16.433	2283	2286	2287	rBV2	11410	10358	0.23%	0.014%
73	16.533	2298	2303	2307	rVB8	16574	25576	0.56%	0.035%
74	16.592	2309	2313	2315	rBV3	27592	41217	0.91%	0.057%
75	16.615	2315	2317	2318	rVV2	22935	20867	0.46%	0.029%

76	16.645	2318	2322	2333	rVB5	61144	124080	2.74%	0.172%
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77	16.851	2351	2357	2363	rVV	2213131	3147502	69.42%	4.367%
78	16.898	2363	2365	2368	rVV	59410	66245	1.46%	0.092%
79	16.951	2368	2374	2384	rVB	3172207	4268703	94.14%	5.923%
80	17.080	2390	2396	2407	rBV6	102883	225978	4.98%	0.314%
81	17.227	2418	2421	2422	rBV3	11080	10773	0.24%	0.015%
82	17.280	2427	2430	2433	rVB4	15746	20157	0.44%	0.028%
83	17.368	2443	2445	2449	rBV4	14689	14514	0.32%	0.020%
84	17.410	2449	2452	2454	rBV4	9325	11934	0.26%	0.017%
85	17.551	2472	2476	2479	rBV2	33384	41583	0.92%	0.058%
86	17.621	2484	2488	2494	rBV6	60023	108551	2.39%	0.151%
87	17.786	2511	2516	2530	rBV2	265874	528102	11.65%	0.733%
88	18.062	2559	2563	2565	rBV2	26088	37192	0.82%	0.052%
89	18.698	2668	2671	2679	rVB	91187	103073	2.27%	0.143%
90	18.915	2705	2708	2711	rBV	87743	99273	2.19%	0.138%
91	19.068	2731	2734	2742	rBV2	74966	136209	3.00%	0.189%
92	19.251	2759	2765	2776	rBV	3381291	4448632	98.11%	6.173%
93	19.815	2858	2861	2866	rVB	442471	474880	10.47%	0.659%
94	20.780	3021	3025	3037	rBV	767827	1113755	24.56%	1.545%
95	21.062	3068	3073	3079	rBV	2025116	2705045	59.66%	3.753%
96	21.768	3189	3193	3195	rBV	310434	427706	9.43%	0.593%
97	21.797	3195	3198	3208	rVB	533827	926010	20.42%	1.285%
98	23.003	3399	3403	3410	rVB	365014	656312	14.47%	0.911%
99	23.597	3497	3504	3512	rBV2	1700153	3721767	82.08%	5.164%
100	23.786	3529	3536	3545	rVB	1316204	2957394	65.22%	4.104%

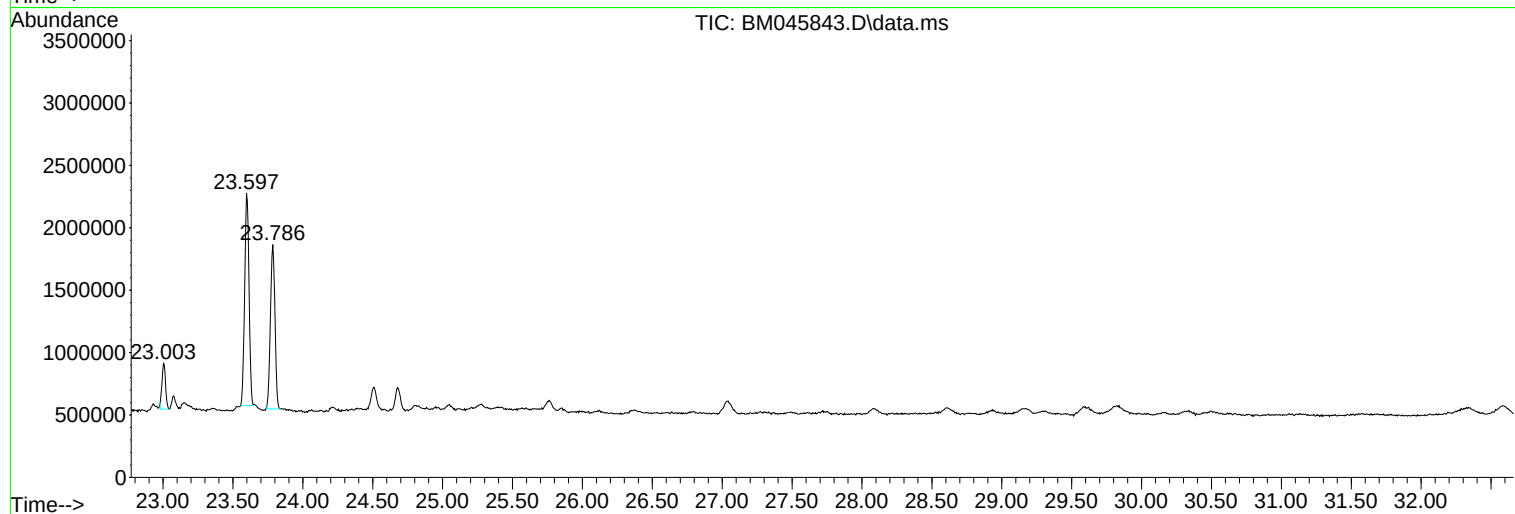
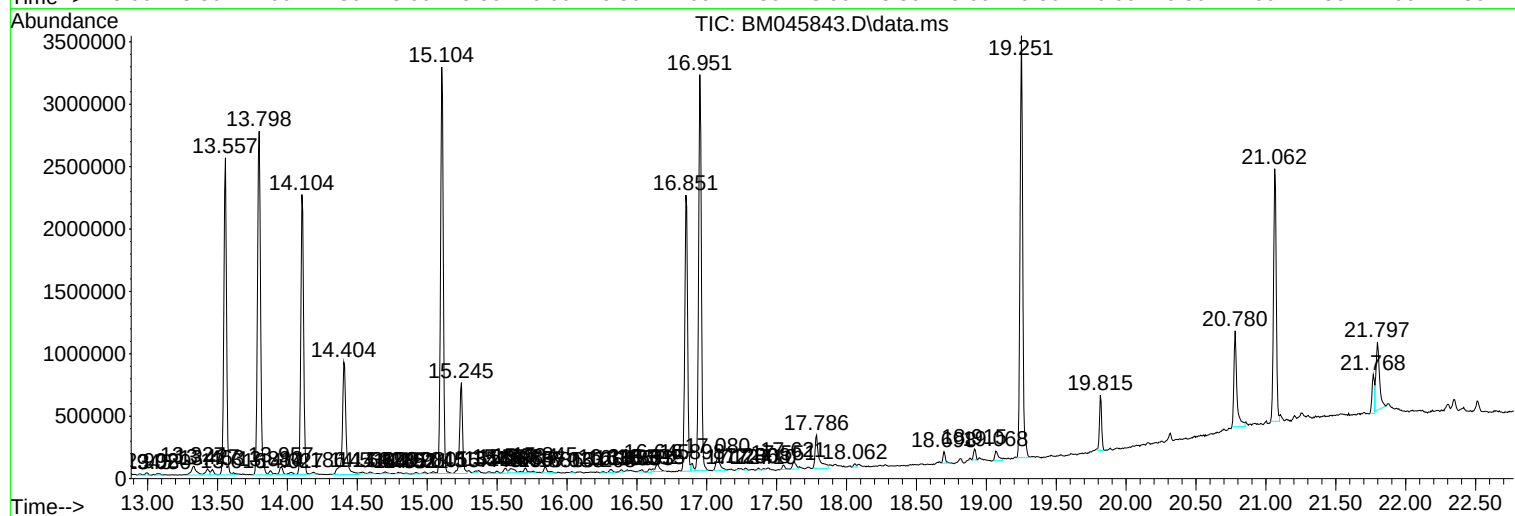
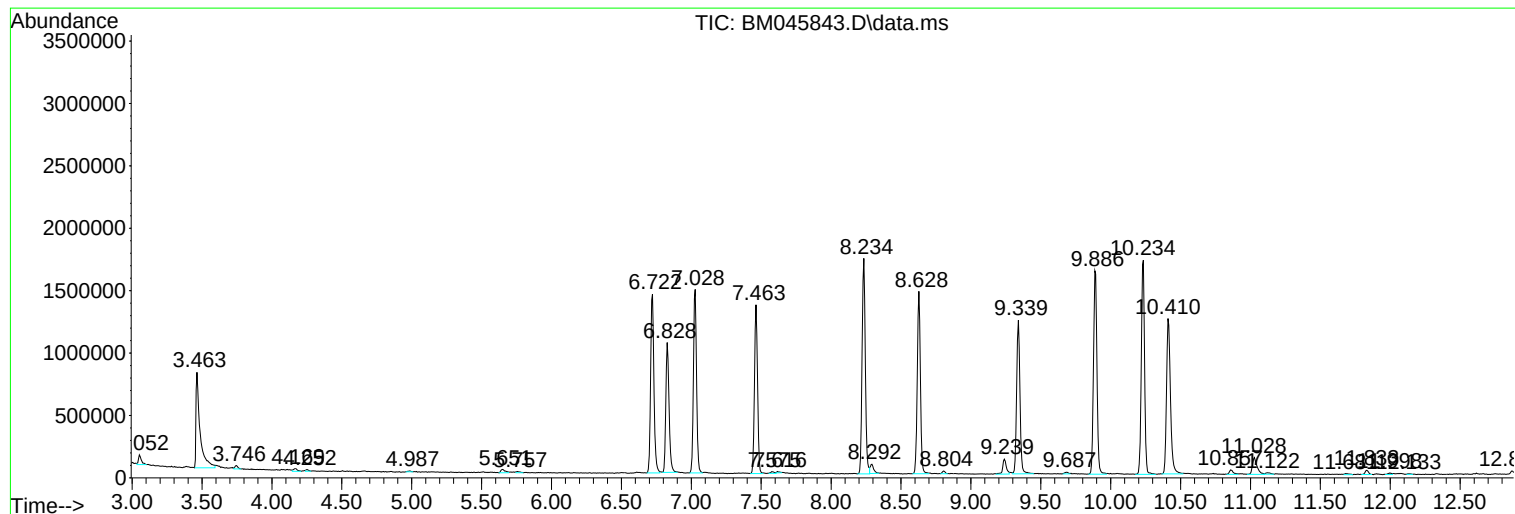
Sum of corrected areas: 72069149

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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



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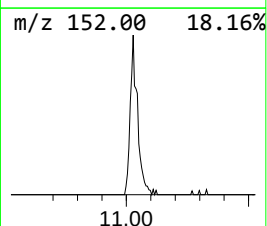
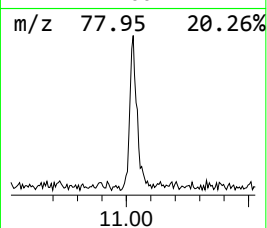
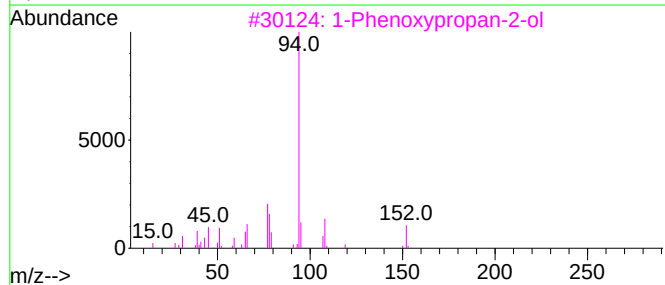
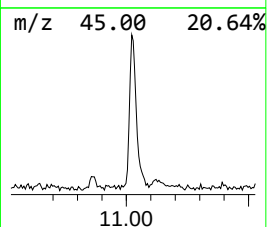
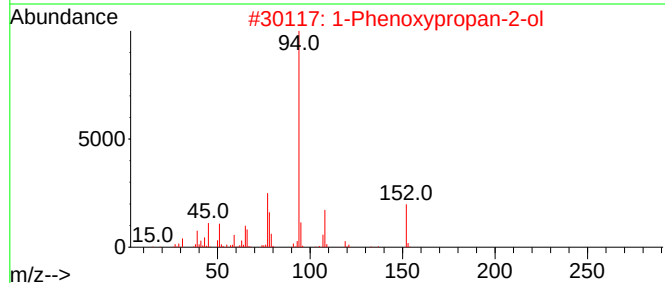
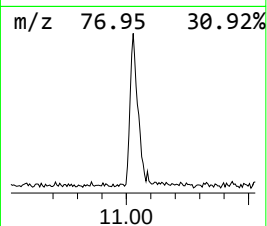
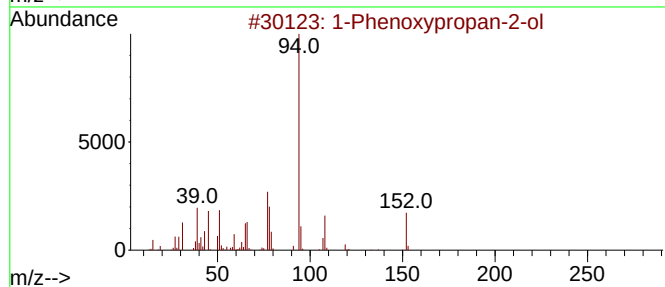
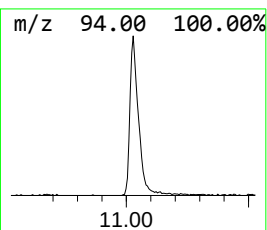
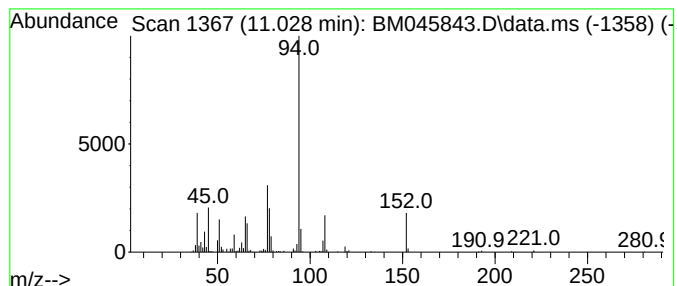
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.028	2.19 ng/ul	308487	Naphthalene-d8	10.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	97
2	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	97
3	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	94
4	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	81
5	1-Propanol, 3-phenoxy-			152	C9H12O2	006180-61-6	64



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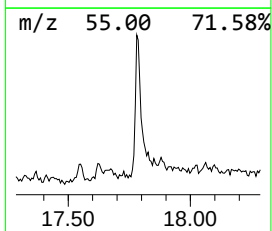
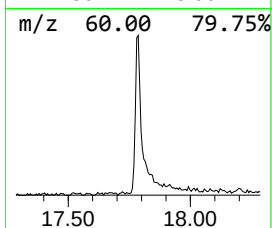
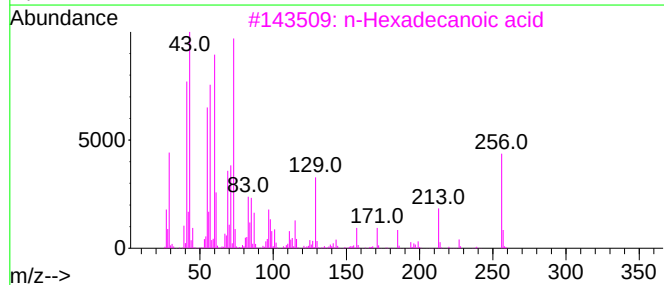
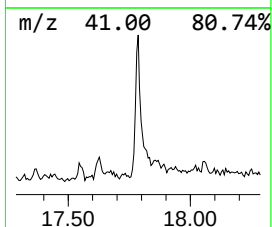
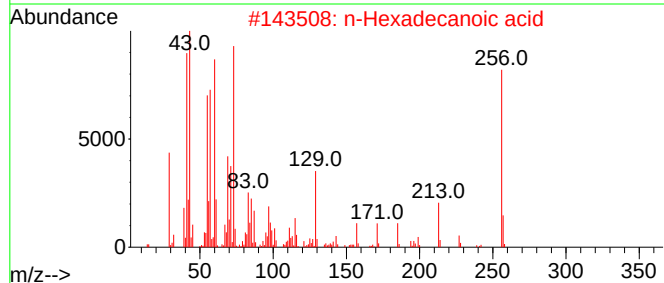
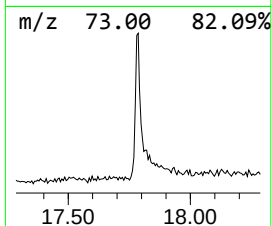
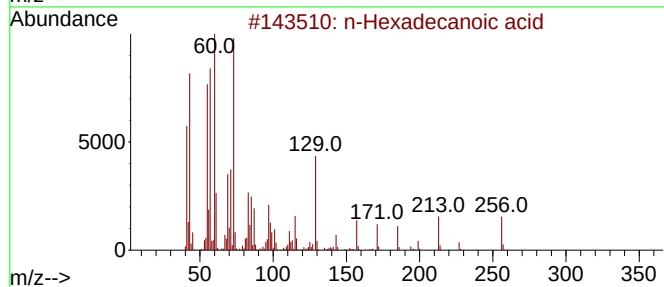
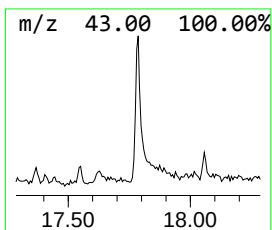
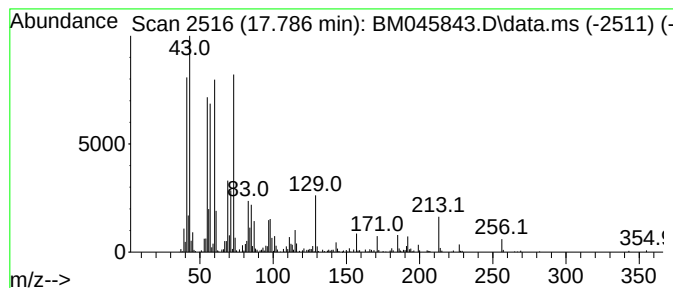
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.786	3.36 ng/ul	528102	Phenanthrene-d10	16.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90		
5	Octadecanoic acid	284	C18H36O2	000057-11-4	90		



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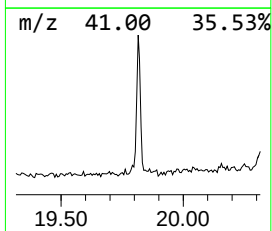
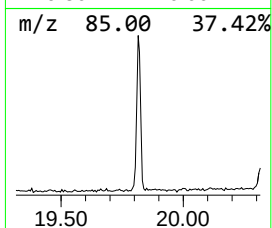
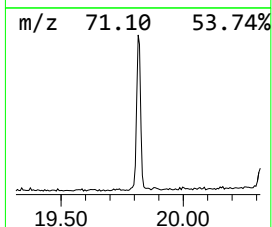
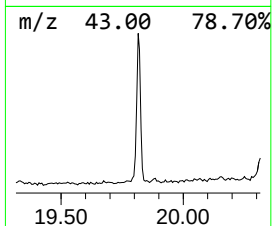
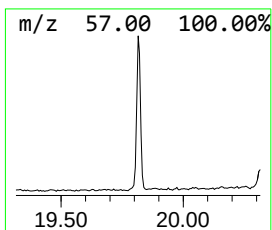
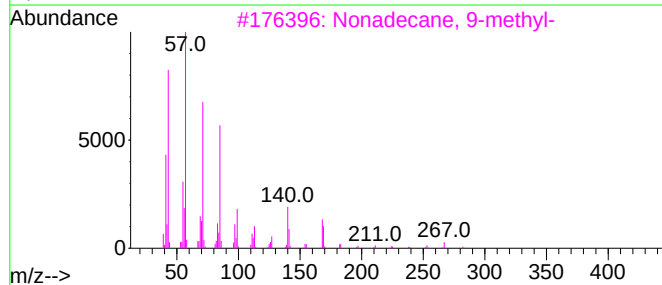
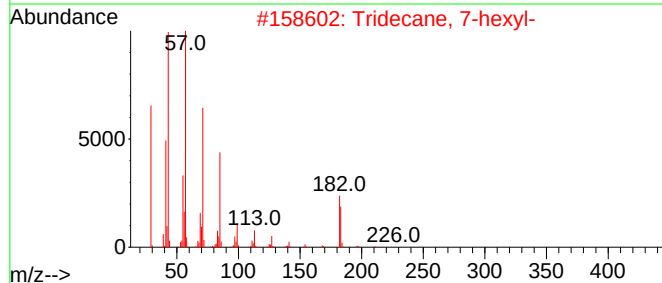
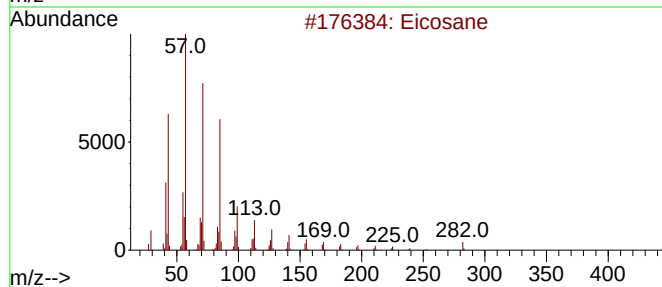
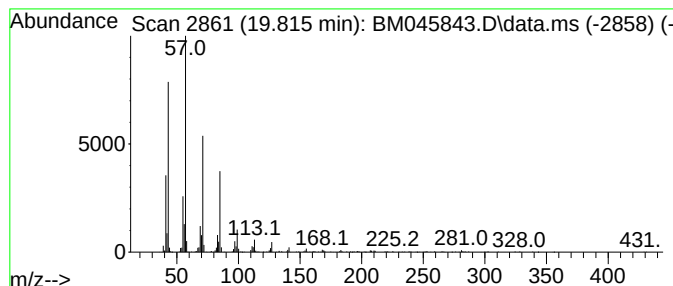
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.815	3.51 ng/ul	474880	Chrysene-d12		21.062	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	Tridecane, 7-hexyl-		268	C19H40	007225-66-3	93
3	Nonadecane, 9-methyl-		282	C20H42	013287-24-6	93
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	93
5	Tridecane, 7-propyl-		226	C16H34	055045-09-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045843.D
Acq On : 05 Jun 2024 09:00
Operator : MA/JU
Sample : P2586-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4D16

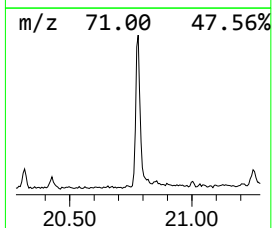
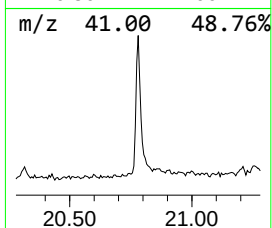
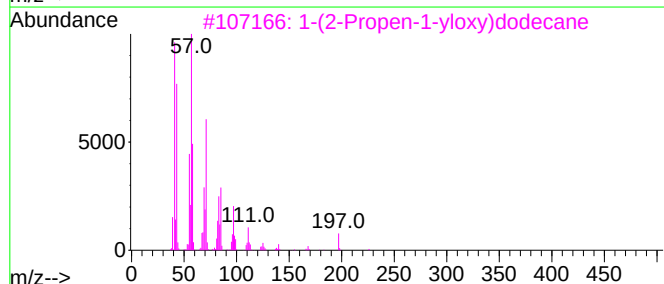
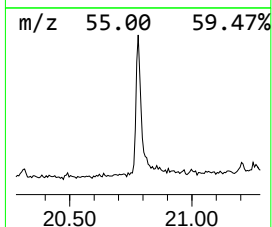
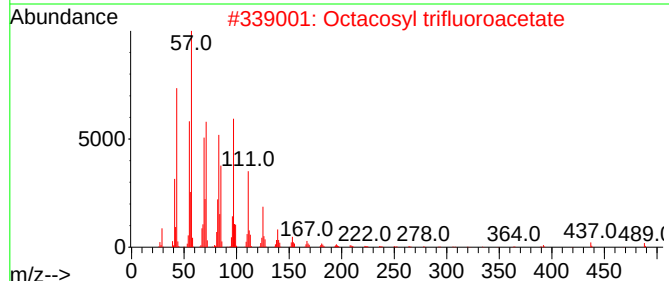
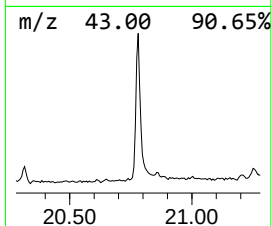
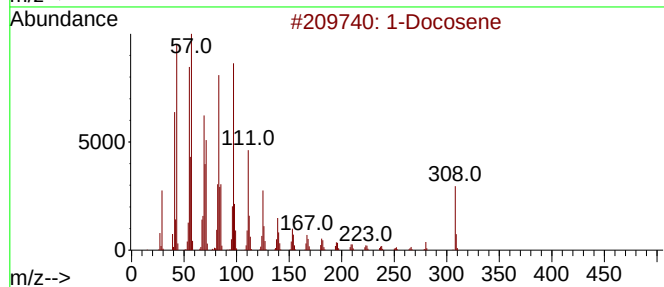
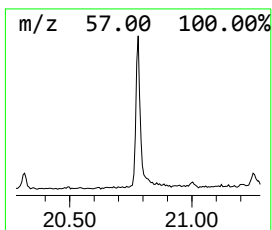
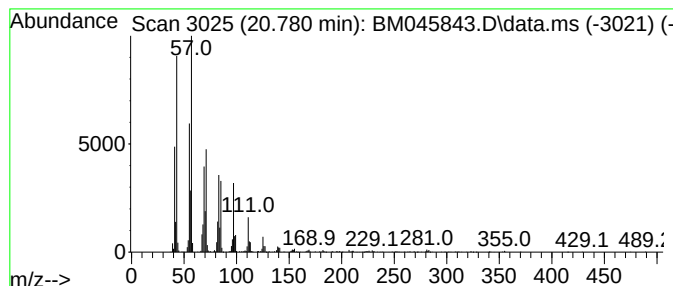
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.780	8.23 ng/ul	1113760	Chrysene-d12	21.062

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene			308	C22H44	001599-67-3	93
2	Octacosyl trifluoroacetate			506	C30H57F3O2	1000351-74-9	91
3	1-(2-Propen-1-yloxy)dodecane			226	C15H30O	006145-80-8	89
4	Oxalic acid, hexadecyl propyl ester			356	C21H40O4	1000309-26-9	87
5	Oxalic acid, pentadecyl propyl e...			342	C20H38O4	1000309-26-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045843.D
Acq On : 05 Jun 2024 09:00
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Sample : P2586-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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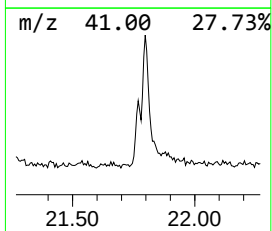
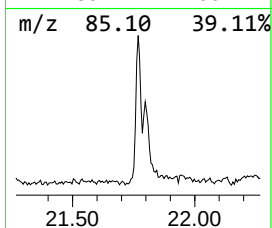
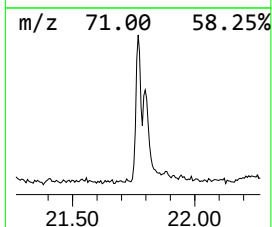
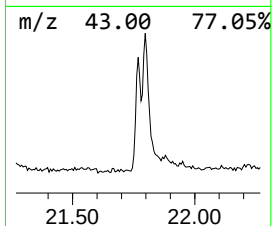
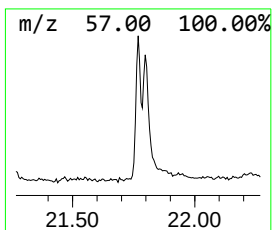
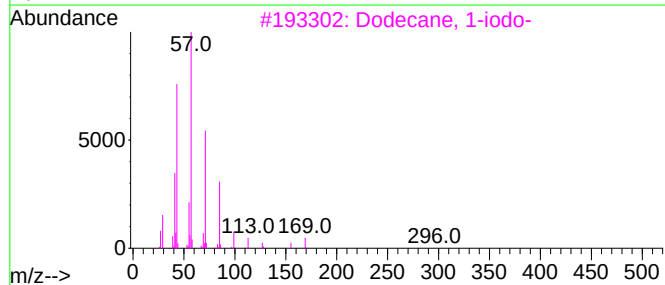
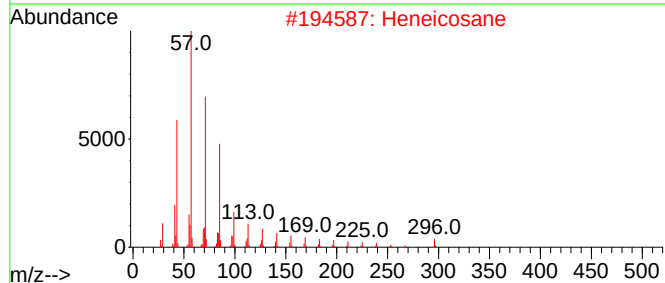
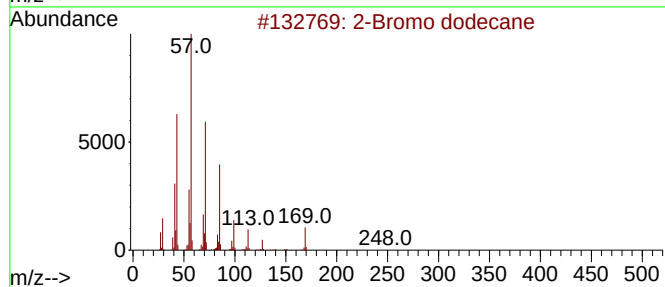
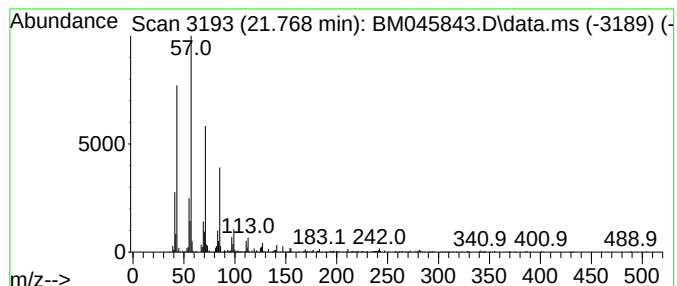
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Bromo dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.768	3.16 ng/ul	427706	Chrysene-d12	21.062

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane			248	C12H25Br	013187-99-0	95
2	Heneicosane			296	C21H44	000629-94-7	95
3	Dodecane, 1-iodo-			296	C12H25I	004292-19-7	93
4	Tridecane, 1-iodo-			310	C13H27I	035599-77-0	93
5	Octacosane			394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045843.D
Acq On : 05 Jun 2024 09:00
Operator : MA/JU
Sample : P2586-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

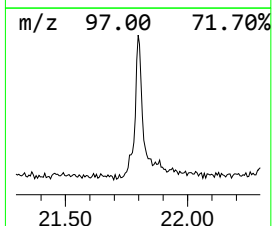
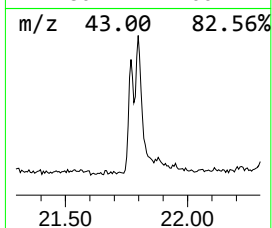
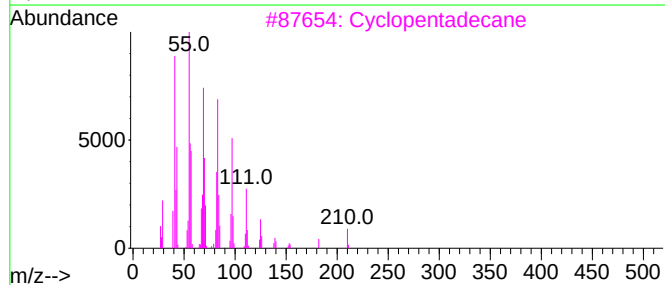
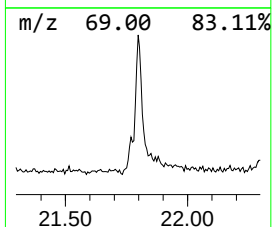
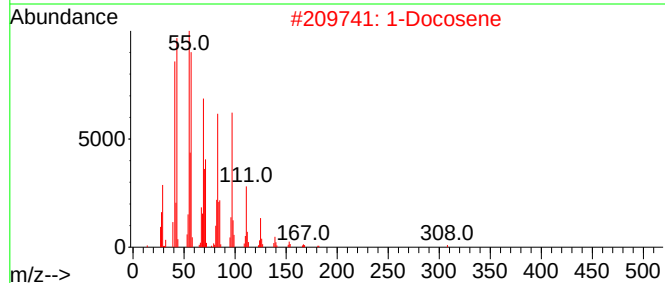
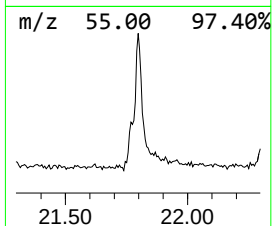
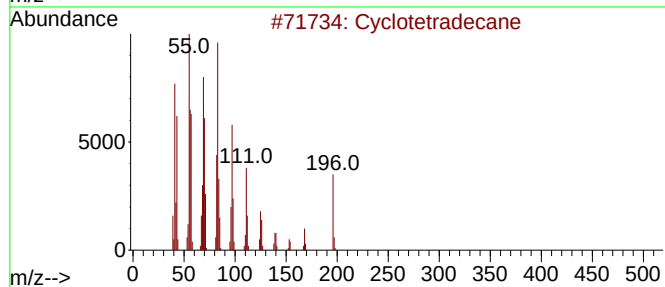
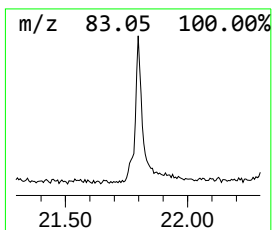
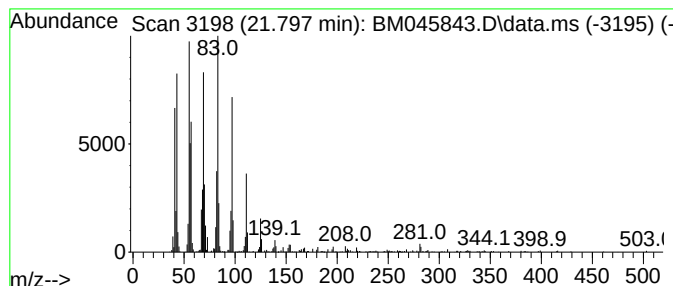
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.797 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.797	6.85 ng/ul	926010	Chrysene-d12	21.062	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	90
2	1-Docosene	308	C22H44	001599-67-3	89
3	Cyclopentadecane	210	C15H30	000295-48-7	87
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	87
5	Z-5-Nonadecene	266	C19H38	1000131-11-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045843.D
Acq On : 05 Jun 2024 09:00
Operator : MA/JU
Sample : P2586-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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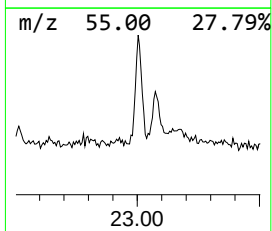
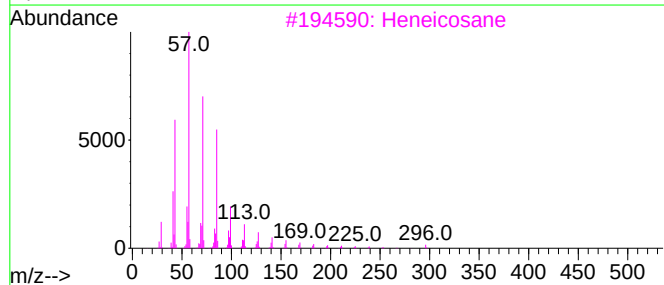
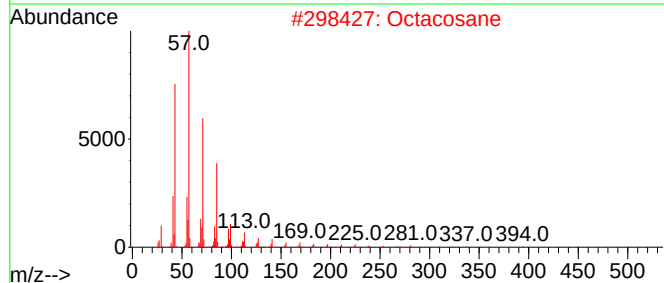
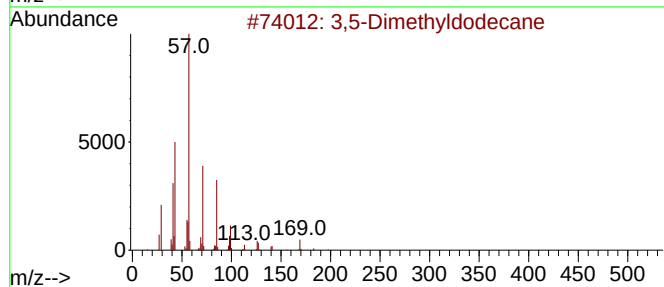
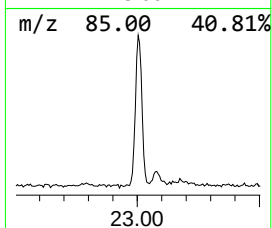
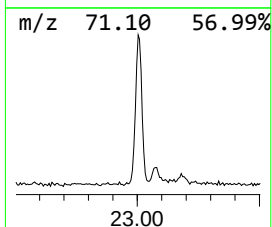
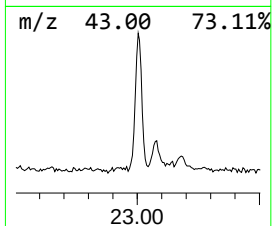
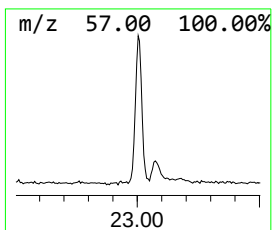
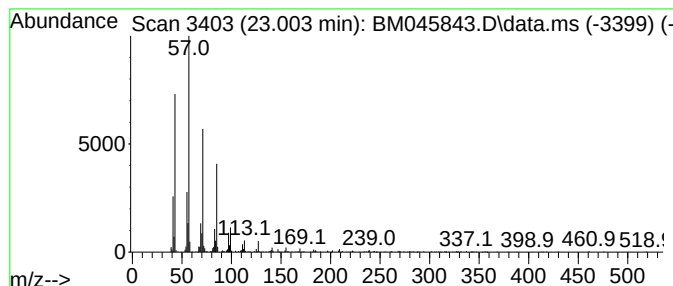
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.003	4.44 ng/ul	656312	Perylene-d12	23.786

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,5-Dimethyldodecane	198	C14H30	107770-99-0	90
2		Octacosane	394	C28H58	000630-02-4	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
Data File : BM045843.D
Acq On : 05 Jun 2024 09:00
Operator : MA/JU
Sample : P2586-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic ...	17.786	3.4	ng/ul	528102	4	16.851	3147500	20.0
(DEL) Alkane: S...	19.815	3.5	ng/ul	474880	5	21.062	2705050	20.0
(DEL) Alkane: S...	20.780	8.2	ng/ul	1113760	5	21.062	2705050	20.0
2-Bromo dodecane	21.768	3.2	ng/ul	427706	5	21.062	2705050	20.0
(DEL) Alkane: C...	21.797	6.8	ng/ul	926010	5	21.062	2705050	20.0
(DEL) Alkane: S...	23.003	4.4	ng/ul	656312	6	23.786	2957390	20.0