

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047021.D
 Acq On : 06 Aug 2024 11:35
 Operator : RC/JU
 Sample : P3171-09
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVX1

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-BM080224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047021.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	19	22	28	rVB	47075	58937	1.44%	0.106%
2	3.281	63	67	73	rBV	95913	141601	3.46%	0.254%
3	3.405	85	88	103	rBV	468684	703754	17.18%	1.261%
4	3.616	119	124	129	rBV	34554	49120	1.20%	0.088%
5	3.705	137	139	145	rVB2	15427	23380	0.57%	0.042%
6	4.287	234	238	241	rBV5	8506	11349	0.28%	0.020%
7	4.599	285	291	297	rBV	207521	289572	7.07%	0.519%
8	5.534	446	450	455	rBV	27774	41187	1.01%	0.074%
9	5.775	485	491	501	rBV	730821	1054710	25.74%	1.890%
10	5.946	516	520	522	rBV5	9125	10189	0.25%	0.018%
11	6.446	599	605	610	rBV6	10495	19598	0.48%	0.035%
12	6.599	624	631	642	rBV	799635	1221965	29.82%	2.189%
13	6.752	650	657	670	rBV	631614	960098	23.43%	1.720%
14	6.934	681	688	698	rBV	796961	1231709	30.06%	2.207%
15	7.028	700	704	715	rVV	48378	86122	2.10%	0.154%
16	7.393	755	766	775	rBV	855334	1300488	31.74%	2.330%
17	7.475	775	780	791	rVB2	61307	110263	2.69%	0.198%
18	7.628	800	806	815	rBV3	65619	128376	3.13%	0.230%
19	7.822	835	839	844	rBV2	22351	33673	0.82%	0.060%
20	8.034	870	875	880	rVB7	7902	16427	0.40%	0.029%
21	8.110	880	888	899	rBV	956267	1528348	37.30%	2.738%
22	8.204	902	904	911	rVV7	7179	11416	0.28%	0.020%
23	8.287	913	918	924	rBV6	11064	19101	0.47%	0.034%
24	8.393	932	936	941	rBV6	6131	9526	0.23%	0.017%
25	8.540	953	961	970	rBV	739818	1207207	29.46%	2.163%
26	8.704	980	989	1000	rBV	96601	173194	4.23%	0.310%
27	8.822	1002	1009	1012	rVV6	11338	22789	0.56%	0.041%
28	8.981	1030	1036	1042	rVB7	9367	20503	0.50%	0.037%
29	9.122	1047	1060	1067	rBV	100102	168404	4.11%	0.302%
30	9.251	1072	1082	1093	rVV	642508	1081036	26.38%	1.937%
31	9.340	1093	1097	1101	rVV	19318	33960	0.83%	0.061%
32	9.404	1101	1108	1117	rVB2	31846	62154	1.52%	0.111%
33	9.510	1117	1126	1137	rBV2	293093	492134	12.01%	0.882%
34	9.687	1151	1156	1164	rBV2	25846	53243	1.30%	0.095%
35	9.787	1166	1173	1181	rVV	991399	1608030	39.25%	2.881%
36	10.057	1214	1219	1225	rBV	68106	107150	2.62%	0.192%

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047021.D
 Acq On : 06 Aug 2024 11:35
 Operator : RC/JU
 Sample : P3171-09
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVX1

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-BM080224.MA.M
 Title : SVOA CALIBRATION

37	10.151	1227	1235	1242	rVV	1075487	1725567	42.11%	3.091%
38	10.222	1242	1247	1253	rVV	72113	114029	2.78%	0.204%
39	10.298	1253	1260	1271	rVV	889282	1467985	35.83%	2.630%
40	10.387	1271	1275	1279	rVB7	9998	17642	0.43%	0.032%

41	10.481	1287	1291	1298	rVB3	19277	33502	0.82%	0.060%
42	10.563	1301	1305	1308	rVV6	6380	9640	0.24%	0.017%
43	10.722	1327	1332	1340	rBV4	34740	56947	1.39%	0.102%
44	10.910	1355	1364	1374	rBV	191533	339542	8.29%	0.608%
45	10.987	1374	1377	1382	rVB7	7824	11564	0.28%	0.021%

46	11.298	1416	1430	1437	rVB2	74347	171264	4.18%	0.307%
47	11.522	1465	1468	1474	rVB6	11698	19243	0.47%	0.034%
48	11.722	1498	1502	1503	rBV4	9970	12599	0.31%	0.023%
49	11.751	1503	1507	1513	rVB2	21268	37123	0.91%	0.067%
50	11.922	1531	1536	1541	rVV2	33321	55799	1.36%	0.100%

51	12.122	1564	1570	1578	rBV8	15574	28758	0.70%	0.052%
52	12.392	1611	1616	1621	rVB7	9071	15019	0.37%	0.027%
53	12.457	1621	1627	1635	rVB4	18311	33863	0.83%	0.061%
54	12.734	1670	1674	1677	rBV5	7113	7848	0.19%	0.014%
55	12.769	1677	1680	1682	rVV4	6319	8793	0.21%	0.016%

56	12.804	1682	1686	1694	rVB2	31775	50513	1.23%	0.090%
57	12.898	1698	1702	1708	rVB8	8577	15029	0.37%	0.027%
58	13.016	1718	1722	1726	rBV3	22810	36658	0.89%	0.066%
59	13.092	1733	1735	1743	rVB9	9714	17998	0.44%	0.032%
60	13.257	1757	1763	1768	rVB8	3933	8675	0.21%	0.016%

61	13.333	1768	1776	1779	rBV	46112	72239	1.76%	0.129%
62	13.363	1779	1781	1787	rVB6	12950	20579	0.50%	0.037%
63	13.475	1794	1800	1811	rVB	1686200	2299801	56.13%	4.120%
64	13.575	1811	1817	1822	rBV6	8785	16819	0.41%	0.030%
65	13.739	1837	1845	1854	rBV	1849115	2806615	68.50%	5.028%

66	13.833	1857	1861	1867	rVB5	12332	22319	0.54%	0.040%
67	13.998	1882	1889	1891	rBV6	6242	13145	0.32%	0.024%
68	14.051	1891	1898	1906	rVB	1562555	2291216	55.92%	4.105%
69	14.292	1931	1939	1950	rBV	816749	1336132	32.61%	2.394%
70	14.386	1952	1955	1960	rVB7	11544	20815	0.51%	0.037%

71	14.510	1973	1976	1984	rVB8	7891	13625	0.33%	0.024%
72	14.592	1984	1990	1997	rBV5	12314	34954	0.85%	0.063%
73	14.692	2002	2007	2013	rVB	140309	198509	4.84%	0.356%
74	14.892	2035	2041	2044	rBV5	10147	18446	0.45%	0.033%
75	15.057	2062	2069	2076	rBV	2320223	3350383	81.77%	6.002%

76	15.192	2085	2092	2106	rBV	717512	1014165	24.75%	1.817%
----	--------	------	------	------	-----	--------	---------	--------	--------

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047021.D
 Acq On : 06 Aug 2024 11:35
 Operator : RC/JU
 Sample : P3171-09
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVX1

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-BM080224.MA.M
 Title : SVOA CALIBRATION

77	15.292	2106	2109	2114	rVB4	12456	18435	0.45%	0.033%
78	15.516	2141	2147	2153	rBV9	5621	9930	0.24%	0.018%
79	15.639	2165	2168	2173	rBV7	8907	13832	0.34%	0.025%
80	15.816	2195	2198	2200	rBV3	10719	12492	0.30%	0.022%
81	15.892	2208	2211	2215	rVB4	16554	19936	0.49%	0.036%
82	15.980	2222	2226	2229	rBV6	5862	9256	0.23%	0.017%
83	16.469	2303	2309	2321	rBV	2038278	2759551	67.35%	4.944%
84	16.610	2328	2333	2336	rBV5	9256	14772	0.36%	0.026%
85	16.810	2360	2367	2376	rBV	2004894	2697848	65.84%	4.833%
86	16.910	2377	2384	2395	rVV	2632952	3559312	86.87%	6.377%
87	17.027	2401	2404	2409	rBV7	7406	12584	0.31%	0.023%
88	17.351	2456	2459	2464	rVV7	6282	11110	0.27%	0.020%
89	17.439	2469	2474	2479	rVV	100478	133082	3.25%	0.238%
90	17.527	2485	2489	2494	rVB5	11403	19337	0.47%	0.035%
91	17.745	2521	2526	2534	rBV	230197	366108	8.94%	0.656%
92	18.780	2699	2702	2708	rVB3	38360	47521	1.16%	0.085%
93	18.857	2711	2715	2719	rVB	42259	53935	1.32%	0.097%
94	19.045	2744	2747	2754	rBV4	68754	110981	2.71%	0.199%
95	19.221	2771	2777	2786	rBV	3161019	4097327	100.00%	7.341%
96	19.315	2791	2793	2799	rVB6	20404	27194	0.66%	0.049%
97	20.780	3038	3042	3049	rBV	288309	387728	9.46%	0.695%
98	21.045	3082	3087	3093	rBV	2091639	2796444	68.25%	5.010%
99	23.574	3509	3517	3528	rVB	1773814	4039035	98.58%	7.236%
100	23.756	3541	3548	3555	rBV	1273230	2683452	65.49%	4.808%

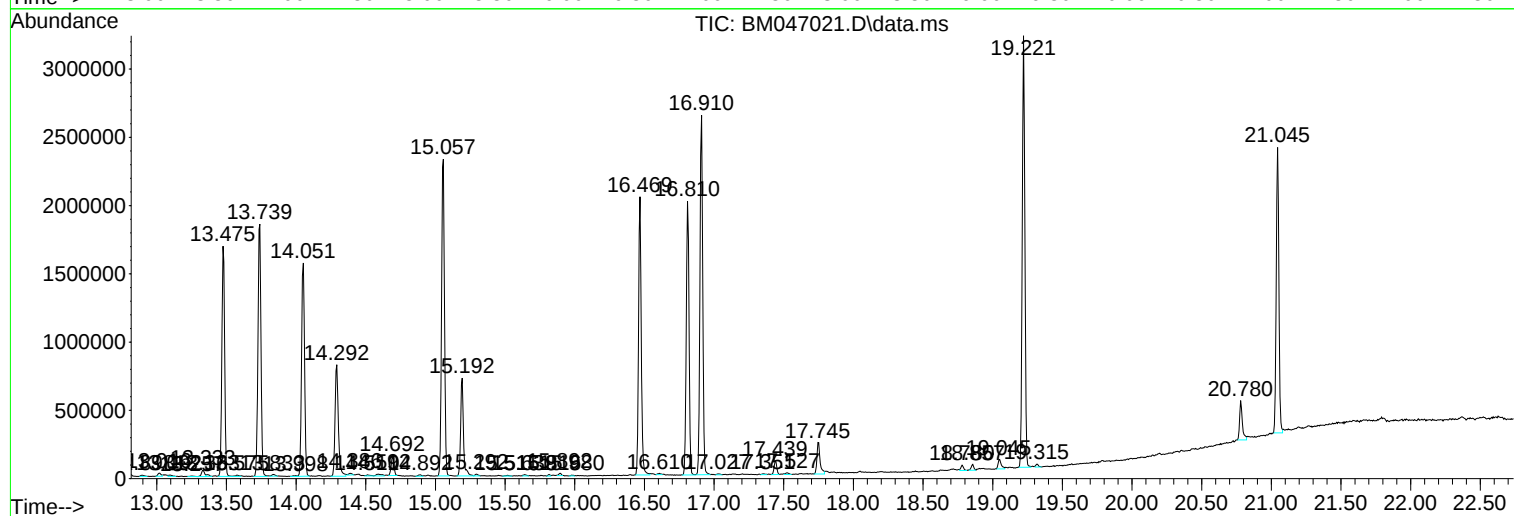
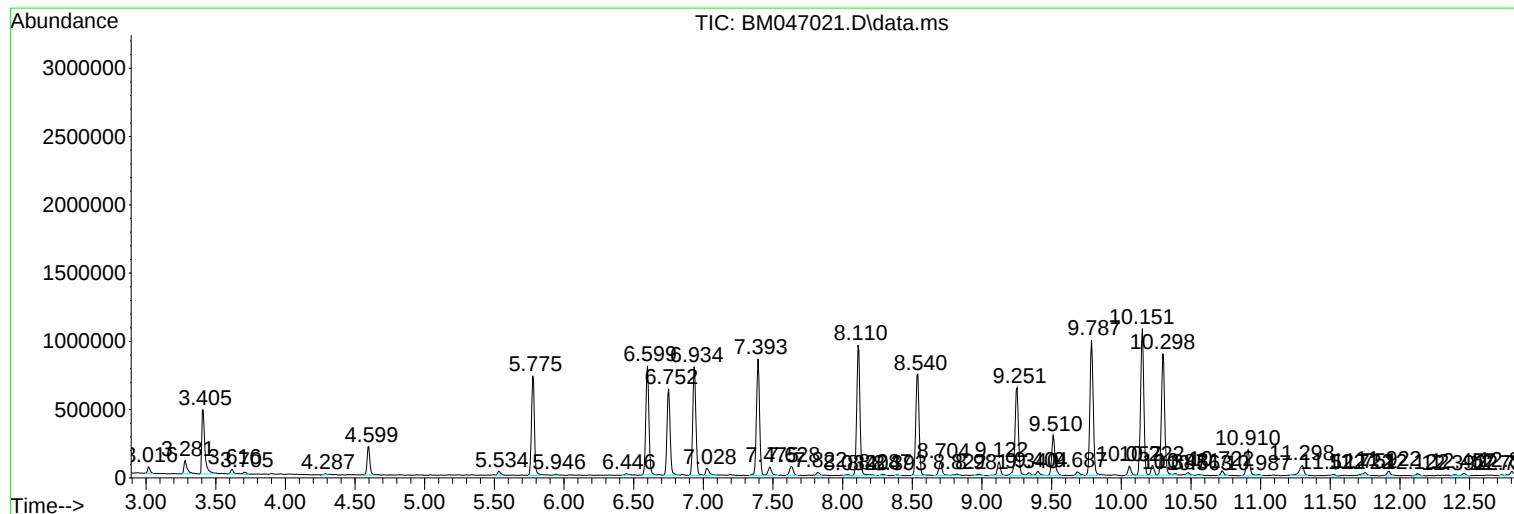
Sum of corrected areas: 55817277

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

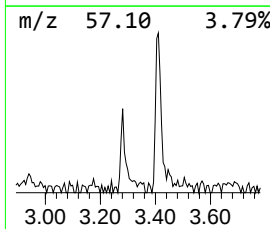
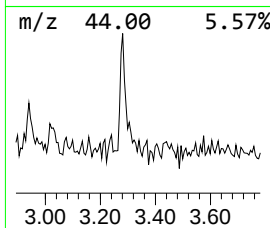
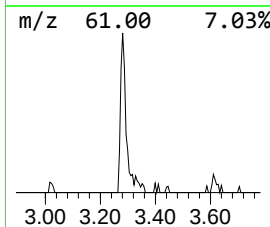
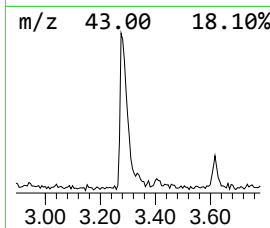
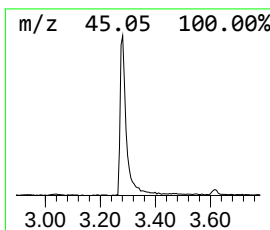
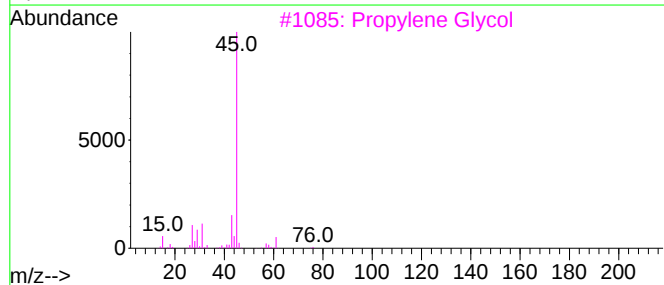
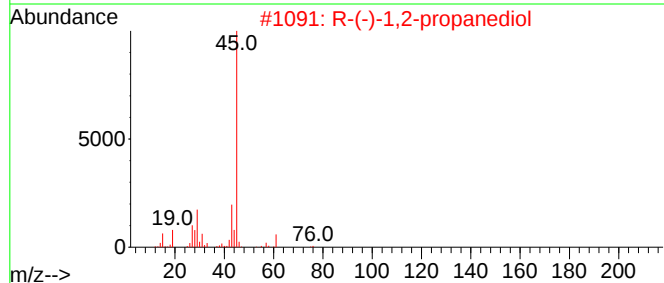
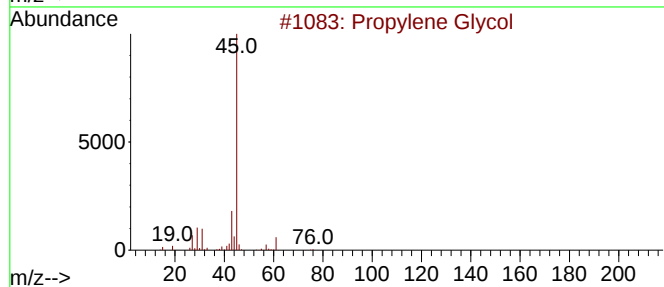
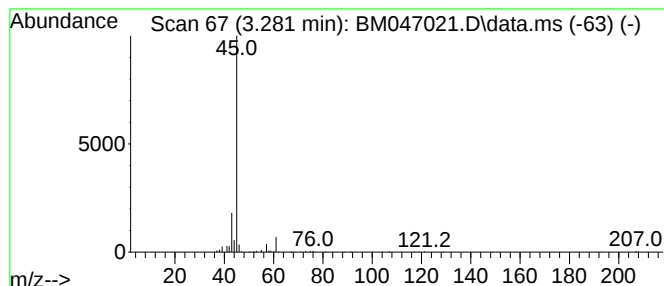
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.281	2.18 ng/ul	141601	1,4-Dichlorobenzene-d4	7.393

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	86
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	72
3			Propylene Glycol	76	C3H8O2	000057-55-6	72
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	50
5			Butanenitrile, 2-(methoxymethoxy)	129	C6H11NO2	1000196-86-5	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

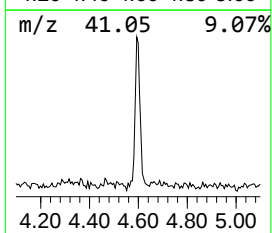
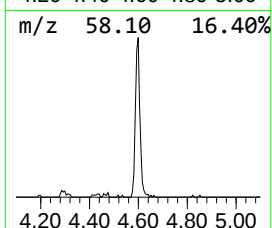
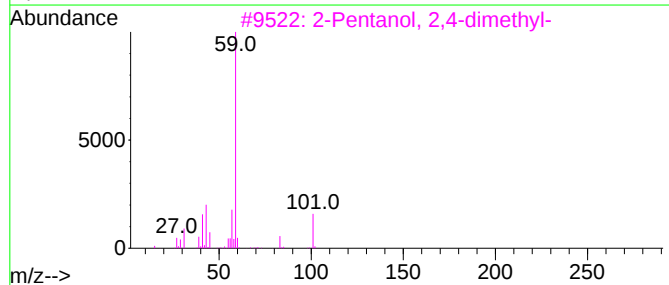
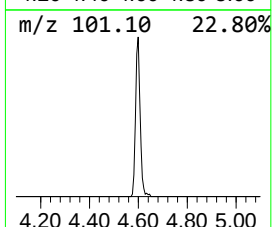
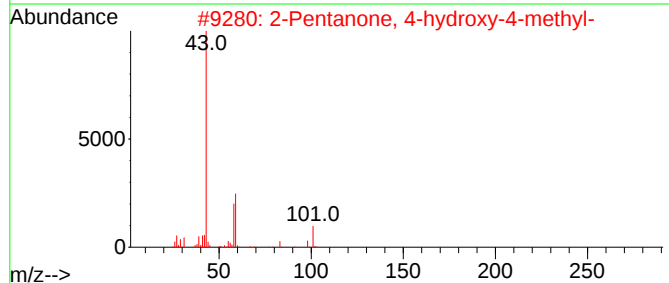
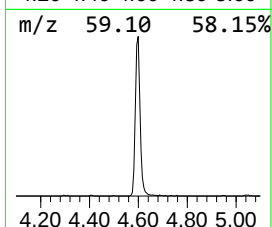
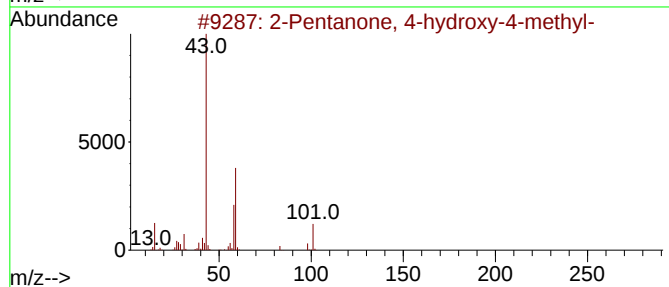
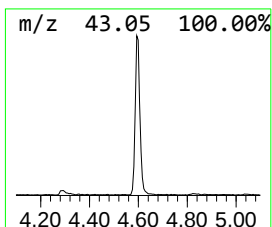
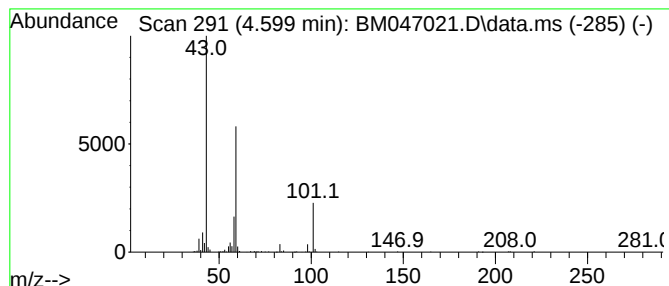
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.599	4.45 ng/ul	289572	1,4-Dichlorobenzene-d4	7.393

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
3	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
5	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

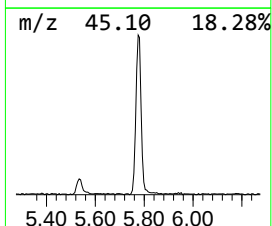
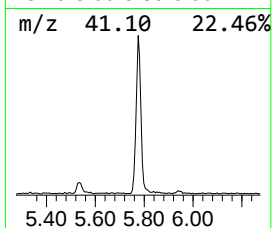
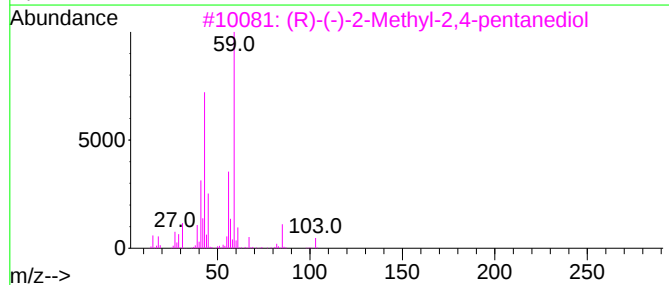
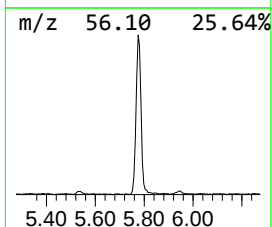
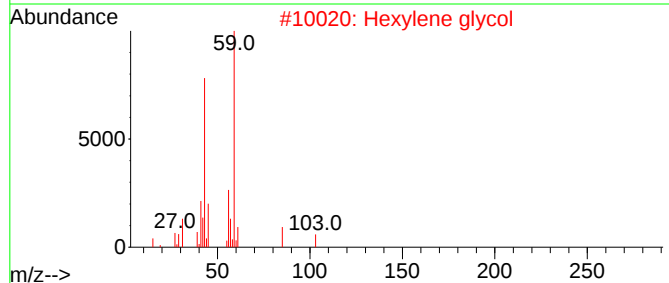
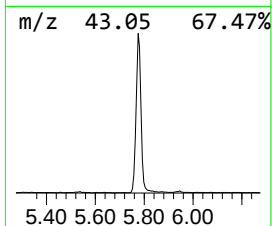
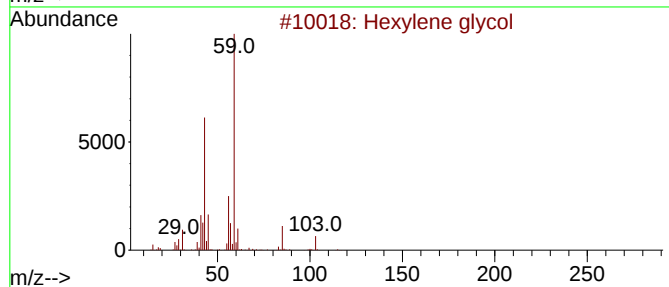
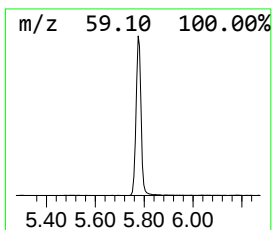
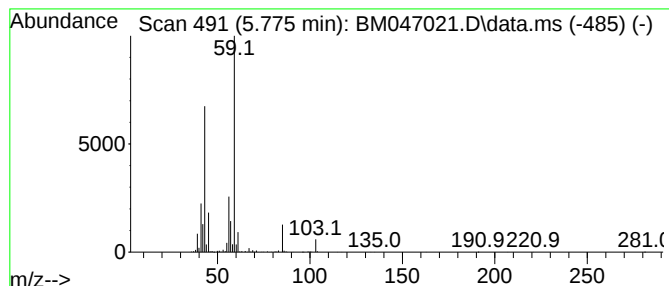
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexylene glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.775	16.22 ng/ul	1054710	1,4-Dichlorobenzene-d4	7.393

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	90
3			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	83
4			Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	78
5			Hexylene glycol	118	C6H14O2	000107-41-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

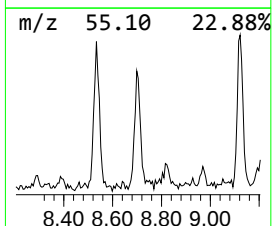
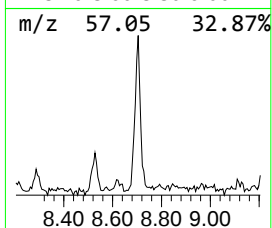
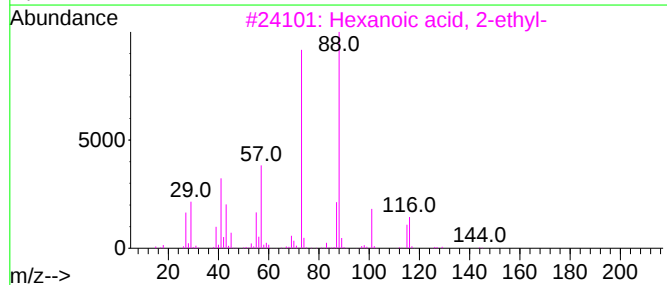
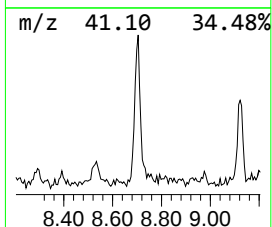
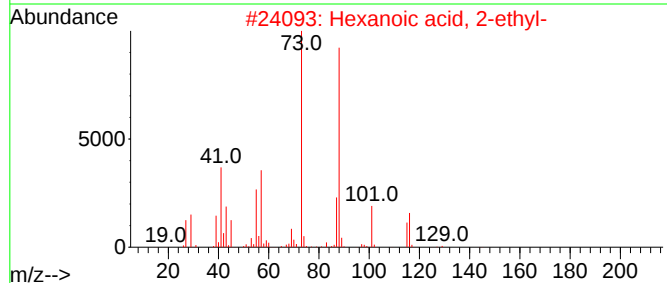
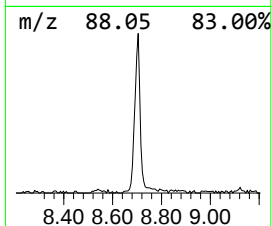
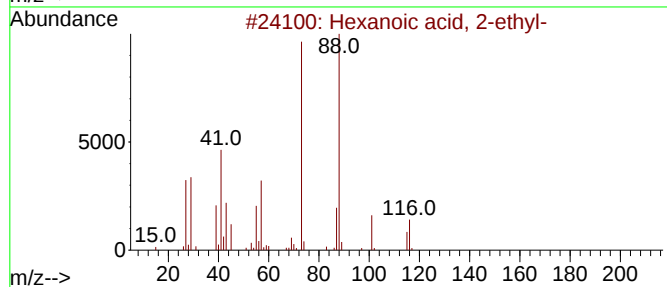
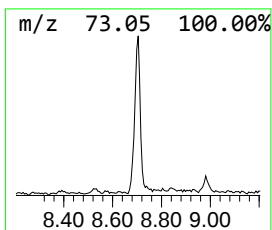
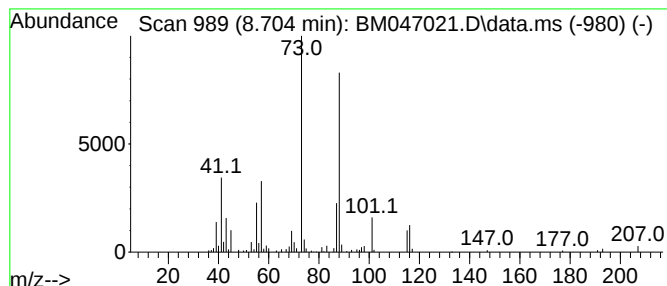
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.704	2.66 ng/ul	173194	1,4-Dichlorobenzene-d4	7.393

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

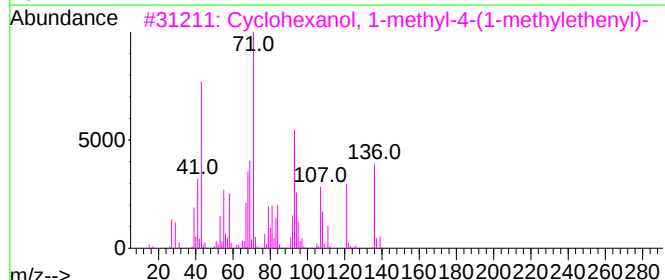
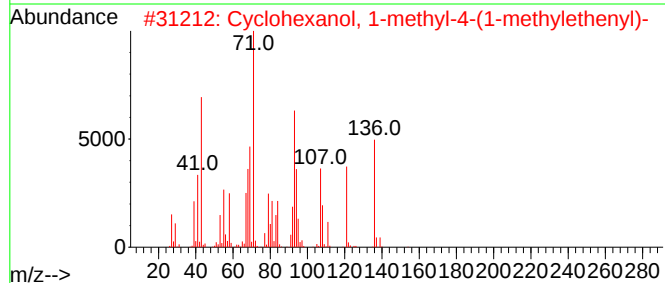
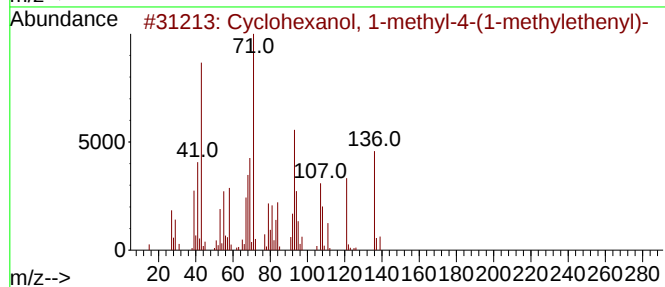
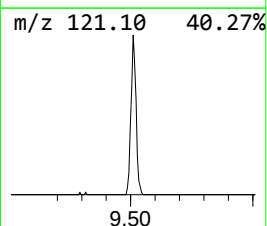
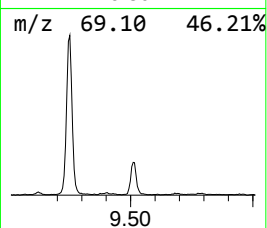
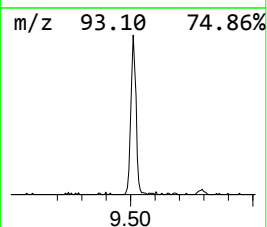
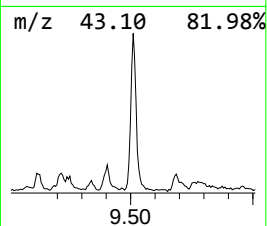
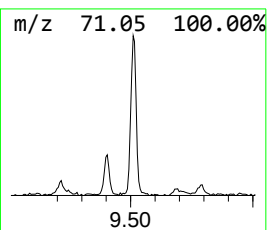
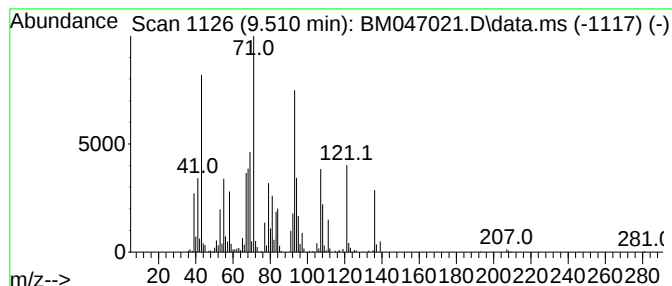
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclohexanol, 1-methyl-4-(1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	5.70 ng/ul	492134	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	90
2			Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	87
3			Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	87
4			Cyclohexene, 5-methyl-3-(1-methy...	136	C10H16	056816-08-1	46
5			Bicyclo[2.2.1]heptane, 7,7-dimet...	136	C10H16	000471-84-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

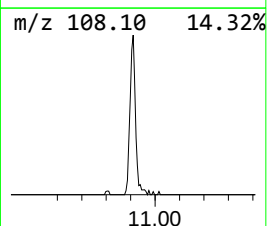
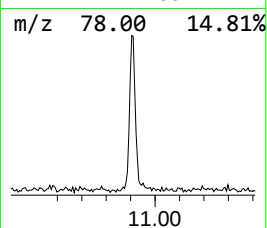
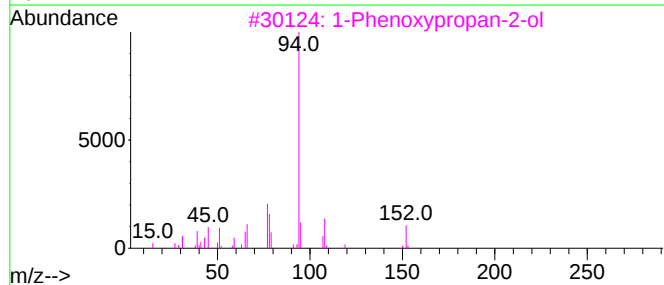
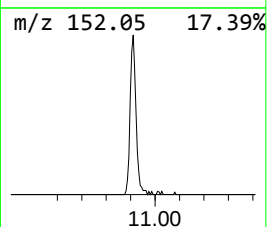
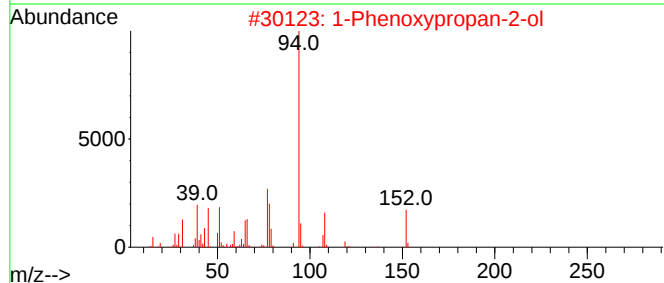
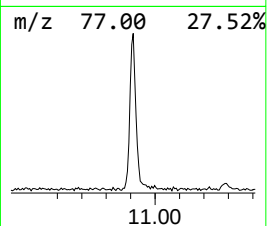
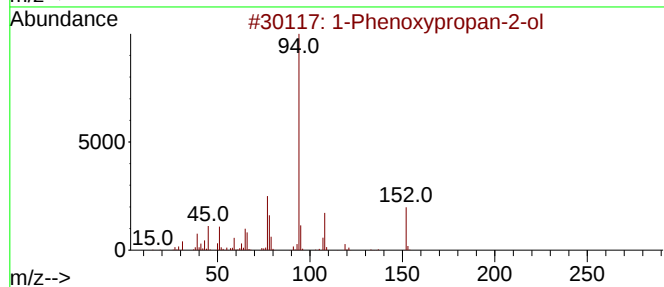
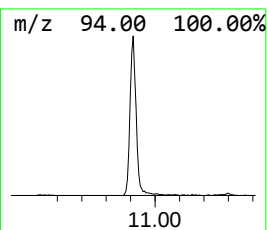
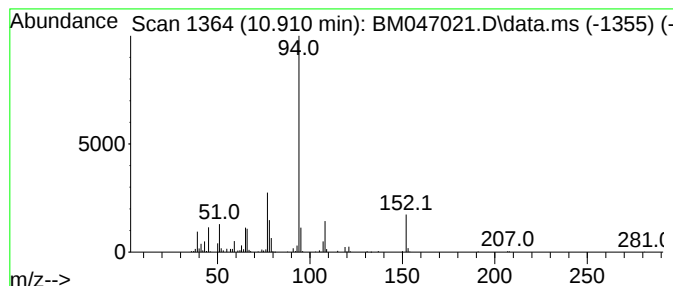
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.910	3.94 ng/ul	339542	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

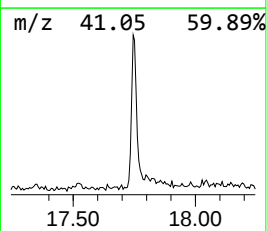
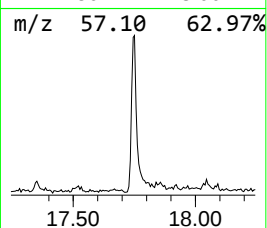
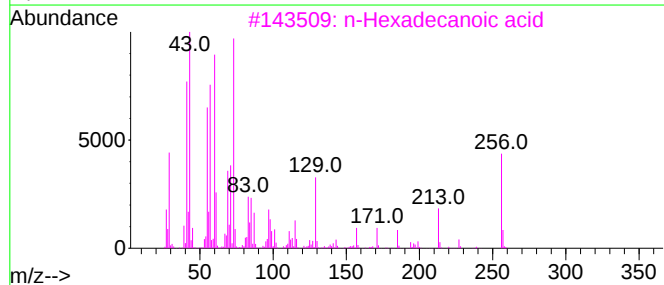
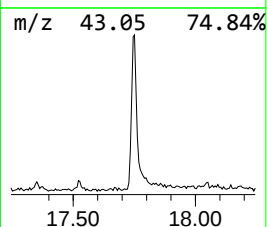
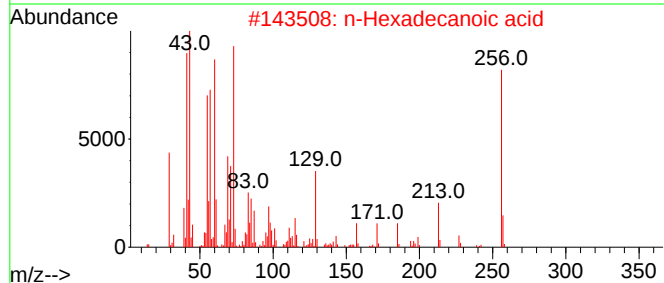
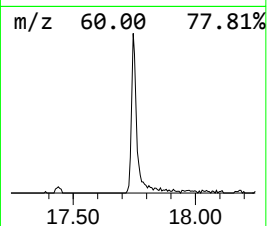
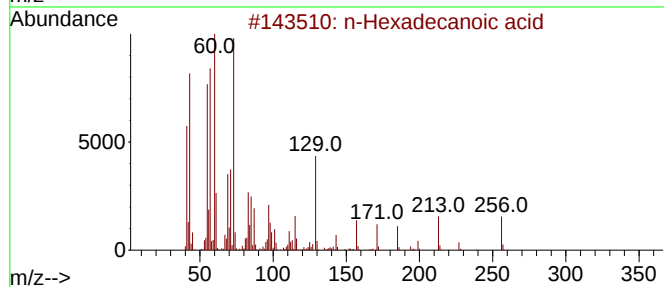
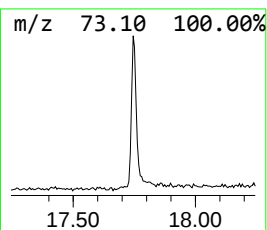
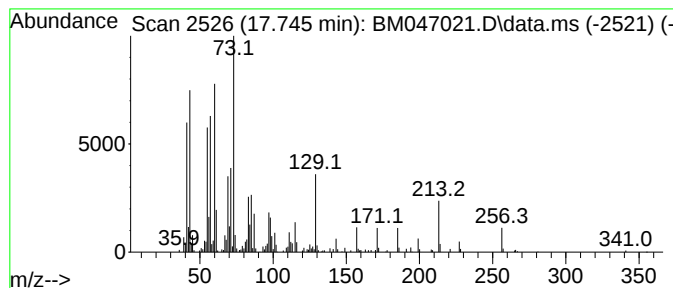
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.745	2.71 ng/ul	366108	Phenanthrene-d10	16.810

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

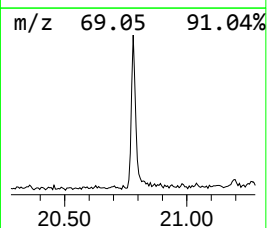
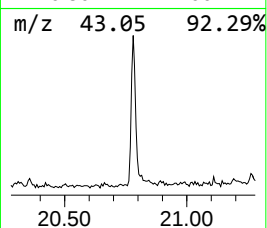
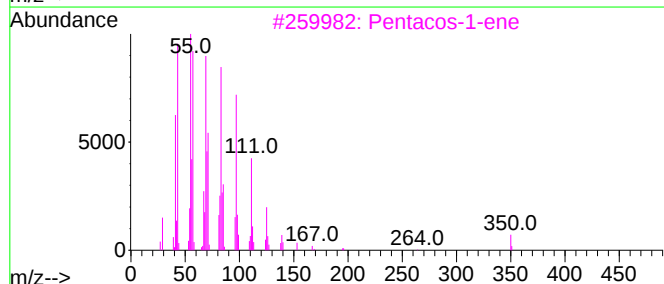
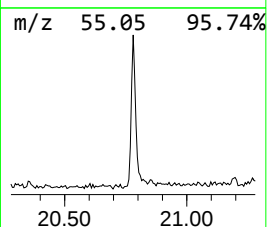
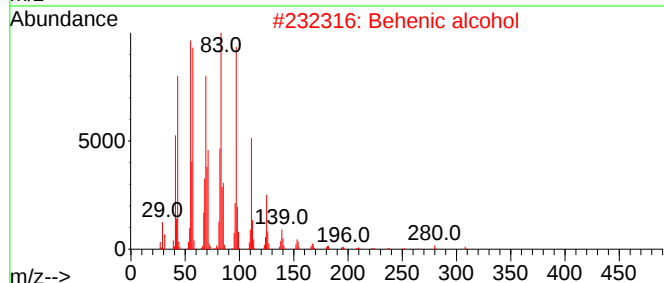
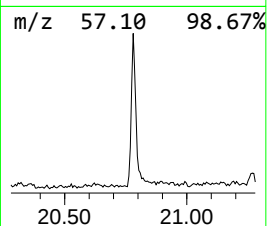
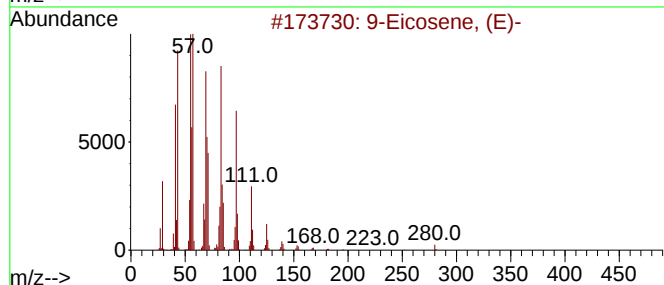
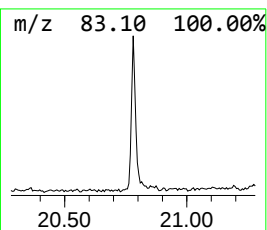
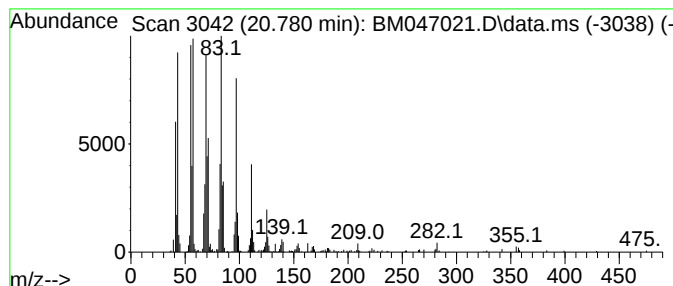
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.780	2.77 ng/ul	387728	Chrysene-d12	21.045

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Eicosene, (E)-		280	C20H40	074685-29-3	95
2	Behenic alcohol		326	C22H46O	000661-19-8	94
3	Pentacos-1-ene		350	C25H50	016980-85-1	94
4	1-Nonadecene		266	C19H38	018435-45-5	92
5	Octacosanol		410	C28H58O	000557-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047021.D
Acq On : 06 Aug 2024 11:35
Operator : RC/JU
Sample : P3171-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.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
Propylene Glycol	3.281	2.2	ng/ul	141601	1	7.393	1300490	20.0
2-Pentanone, 4-...	4.599	4.5	ng/ul	289572	1	7.393	1300490	20.0
Hexylene glycol	5.775	16.2	ng/ul	1054710	1	7.393	1300490	20.0
Hexanoic acid, ...	8.704	2.7	ng/ul	173194	1	7.393	1300490	20.0
Cyclohexanol, 1...	9.510	5.7	ng/ul	492134	2	10.151	1725570	20.0
1-Phenoxypropan...	10.910	3.9	ng/ul	339542	2	10.151	1725570	20.0
n-Hexadecanoic ...	17.745	2.7	ng/ul	366108	4	16.810	2697850	20.0
(DEL) Alkane: S...	20.780	2.8	ng/ul	387728	5	21.045	2796440	20.0