

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
 Data File : BM042754.D
 Acq On : 15 Nov 2023 03:59
 Operator : MA/JU
 Sample : 05079-02
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BHEF1

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-BM111323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM042754.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	30	33	37	rBV3	10939	14185	0.94%	0.104%
2	3.457	76	80	89	rVB	12373	19855	1.31%	0.146%
3	3.628	105	109	122	rBV	55012	96845	6.41%	0.710%
4	4.257	212	216	223	rVB7	4614	9318	0.62%	0.068%
5	4.446	246	248	251	rVB4	2574	2823	0.19%	0.021%
6	4.769	301	303	309	rVB6	2278	3601	0.24%	0.026%
7	4.987	336	340	345	rVB6	4431	6942	0.46%	0.051%
8	5.063	347	353	362	rVB	23518	37862	2.51%	0.278%
9	5.263	383	387	391	rBV3	2798	4593	0.30%	0.034%
10	6.628	615	619	622	rBV5	1545	2656	0.18%	0.019%
11	6.804	645	649	653	rVB5	1824	2486	0.16%	0.018%
12	7.004	677	683	702	rBV	51382	117889	7.80%	0.865%
13	7.181	702	713	726	rBV	153661	280977	18.60%	2.061%
14	7.281	726	730	735	rVB7	2440	3453	0.23%	0.025%
15	7.351	737	742	756	rBV	185559	326214	21.59%	2.392%
16	7.828	817	823	835	rBV	131288	235896	15.61%	1.730%
17	7.934	837	841	849	rVB7	5372	10859	0.72%	0.080%
18	8.186	877	884	889	rBV2	5564	9342	0.62%	0.069%
19	8.310	899	905	907	rBV5	2241	4692	0.31%	0.034%
20	8.492	932	936	937	rBV4	2613	2925	0.19%	0.021%
21	8.557	940	947	959	rBV	129303	264166	17.48%	1.937%
22	8.716	970	974	975	rBV4	3857	3554	0.24%	0.026%
23	8.775	979	984	990	rVB5	10357	16274	1.08%	0.119%
24	8.845	990	996	1003	rVB2	16269	25792	1.71%	0.189%
25	8.963	1011	1016	1022	rVB6	2250	4512	0.30%	0.033%
26	9.033	1022	1028	1033	rBV	197663	359512	23.79%	2.637%
27	9.104	1033	1040	1051	rVB	96040	231622	15.33%	1.699%
28	9.222	1058	1060	1067	rVV7	1466	3436	0.23%	0.025%
29	9.275	1067	1069	1073	rVB5	3263	3748	0.25%	0.027%
30	9.369	1083	1085	1089	rVB5	2522	2844	0.19%	0.021%
31	9.422	1092	1094	1099	rVB6	2484	3205	0.21%	0.024%
32	9.616	1120	1127	1131	rBV8	3947	9687	0.64%	0.071%
33	9.675	1131	1137	1142	rVB6	9214	16755	1.11%	0.123%
34	9.745	1142	1149	1160	rBV	174395	328627	21.75%	2.410%
35	9.916	1175	1178	1179	rBV3	2931	2709	0.18%	0.020%
36	10.169	1215	1221	1223	rBV3	4154	4600	0.30%	0.034%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
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37	10.216	1226	1229	1232	rVB5	2232	2933	0.19%	0.022%
38	10.269	1232	1238	1259	rBV	228309	489000	32.36%	3.586%
39	10.457	1263	1270	1273	rBV5	12567	23326	1.54%	0.171%
40	10.504	1273	1278	1293	rVB3	27402	60003	3.97%	0.440%

41	10.645	1295	1302	1311	rBV	195569	331574	21.94%	2.432%
42	10.763	1318	1322	1324	rBV4	4806	7536	0.50%	0.055%
43	10.827	1327	1333	1341	rBV2	45220	110315	7.30%	0.809%
44	11.022	1359	1366	1369	rBV8	8431	22832	1.51%	0.167%
45	11.351	1414	1422	1431	rBV9	8382	23870	1.58%	0.175%

46	11.451	1431	1439	1441	rBV7	3308	7936	0.53%	0.058%
47	11.533	1449	1453	1456	rVB5	3601	6114	0.40%	0.045%
48	11.604	1461	1465	1468	rBV5	2263	3178	0.21%	0.023%
49	11.774	1490	1494	1499	rBV3	7879	13657	0.90%	0.100%
50	11.816	1499	1501	1506	rVB6	3123	3392	0.22%	0.025%

51	11.916	1510	1518	1526	rVB7	6485	19435	1.29%	0.143%
52	12.021	1528	1536	1541	rBV5	7592	17169	1.14%	0.126%
53	12.086	1542	1547	1548	rBV5	2018	2770	0.18%	0.020%
54	12.186	1560	1564	1565	rBV3	3270	4961	0.33%	0.036%
55	12.304	1579	1584	1591	rVB9	7604	16355	1.08%	0.120%

56	12.369	1591	1595	1600	rVB7	2886	4993	0.33%	0.037%
57	12.445	1604	1608	1609	rBV3	3930	4656	0.31%	0.034%
58	12.510	1614	1619	1623	rBV5	8175	15952	1.06%	0.117%
59	12.551	1623	1626	1630	rVB6	6295	8240	0.55%	0.060%
60	12.674	1641	1647	1655	rBV3	20083	37257	2.47%	0.273%

61	12.792	1663	1667	1671	rBV5	5246	8774	0.58%	0.064%
62	13.027	1702	1707	1710	rBV6	4808	8337	0.55%	0.061%
63	13.063	1711	1713	1720	rVB7	4757	8093	0.54%	0.059%
64	13.210	1733	1738	1740	rBV5	3306	5170	0.34%	0.038%
65	13.257	1743	1746	1751	rBV7	2776	5115	0.34%	0.038%

66	13.327	1754	1758	1765	rVB6	9122	17464	1.16%	0.128%
67	13.416	1771	1773	1777	rVB5	2449	2523	0.17%	0.019%
68	13.468	1777	1782	1783	rBV3	3496	4756	0.31%	0.035%
69	13.504	1785	1788	1792	rVB5	5285	6968	0.46%	0.051%
70	13.663	1812	1815	1817	rBV3	2487	3107	0.21%	0.023%

71	13.745	1827	1829	1832	rVB4	2668	2847	0.19%	0.021%
72	13.798	1832	1838	1842	rBV5	16542	27970	1.85%	0.205%
73	13.910	1851	1857	1863	rBV	586395	759132	50.24%	5.567%
74	14.063	1880	1883	1884	rBV2	3144	3052	0.20%	0.022%
75	14.186	1897	1904	1917	rBV	560943	827545	54.77%	6.069%

76	14.351	1930	1932	1939	rBV7	3024	5156	0.34%	0.038%
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 Title : SVOA CALIBRATION

77	14.486	1949	1955	1963	rBV	300597	443204	29.33%	3.250%
78	14.768	1998	2003	2012	rBV7	13436	44565	2.95%	0.327%
79	14.845	2014	2016	2027	rVV7	8234	19548	1.29%	0.143%
80	14.933	2027	2031	2035	rVV7	3815	7504	0.50%	0.055%
81	14.986	2037	2040	2044	rBV6	5302	7718	0.51%	0.057%
82	15.021	2044	2046	2050	rBV5	4477	6215	0.41%	0.046%
83	15.062	2050	2053	2057	rVB6	4847	8228	0.54%	0.060%
84	15.186	2072	2074	2077	rVV3	7802	11319	0.75%	0.083%
85	15.221	2077	2080	2085	rVB6	8553	13461	0.89%	0.099%
86	15.480	2118	2124	2134	rBV	732995	1029304	68.12%	7.549%
87	15.574	2136	2140	2144	rVV5	10620	20355	1.35%	0.149%
88	15.627	2144	2149	2159	rVV	254002	377615	24.99%	2.769%
89	15.857	2184	2188	2195	rVB	33841	45739	3.03%	0.335%
90	16.015	2213	2215	2216	rBV2	3169	2702	0.18%	0.020%
91	16.409	2279	2282	2287	rVB	16666	22530	1.49%	0.165%
92	17.239	2418	2423	2434	rBV	356092	535233	35.42%	3.925%
93	17.339	2434	2440	2449	rBV	769711	1103460	73.03%	8.092%
94	18.098	2564	2569	2576	rBV	135183	218275	14.45%	1.601%
95	19.374	2783	2786	2795	rBV	61619	95370	6.31%	0.699%
96	19.633	2825	2830	2844	rBV	1093832	1510957	100.00%	11.081%
97	21.086	3074	3077	3083	rVB2	93342	103963	6.88%	0.762%
98	21.421	3130	3134	3146	rBV	441084	566566	37.50%	4.155%
99	23.633	3503	3510	3522	rBV	699377	1392976	92.19%	10.216%
100	23.785	3530	3536	3547	rVB	314806	641045	42.43%	4.701%

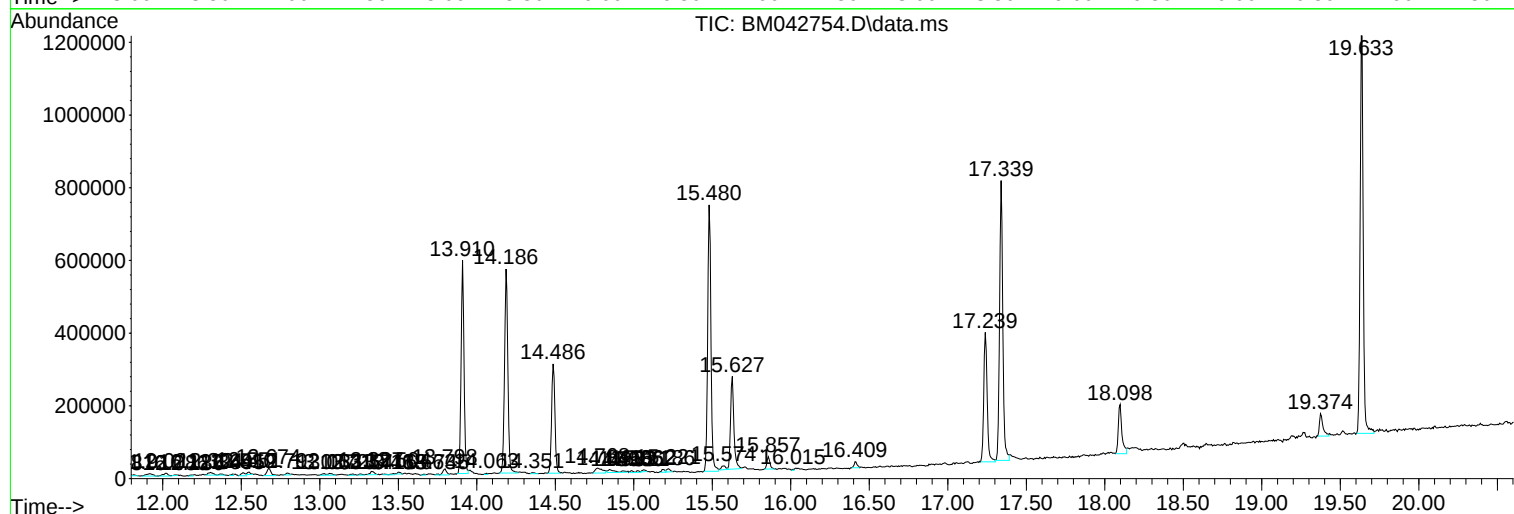
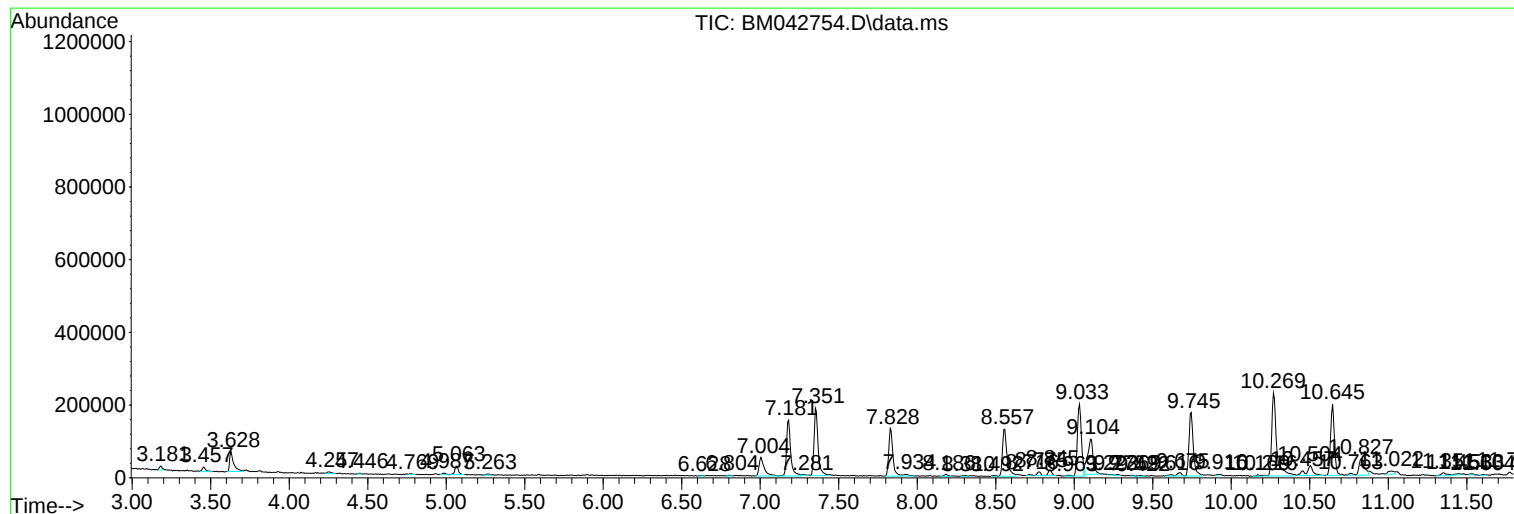
Sum of corrected areas: 13635771

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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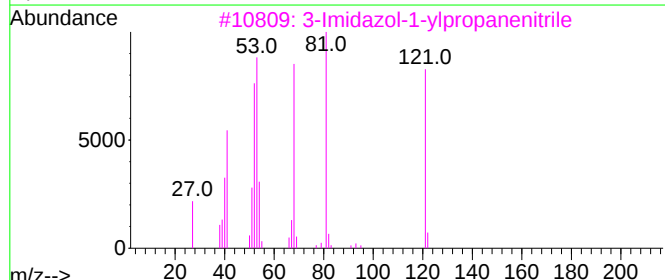
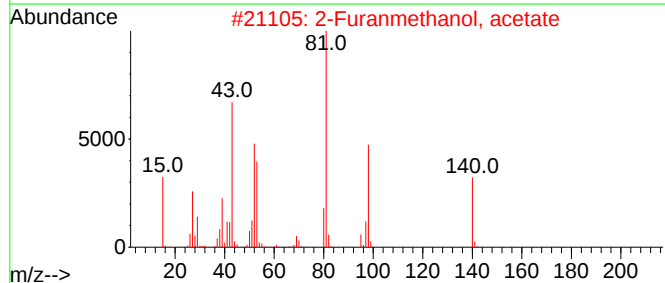
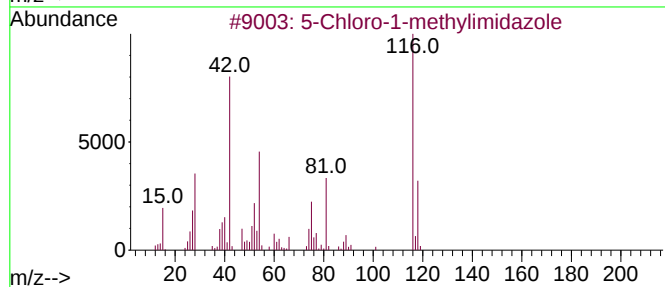
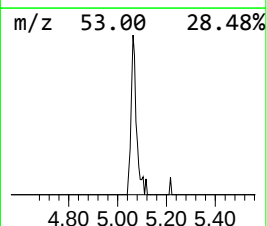
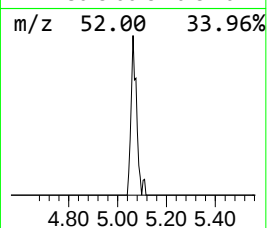
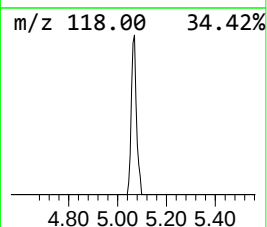
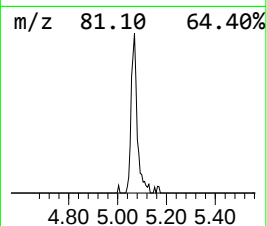
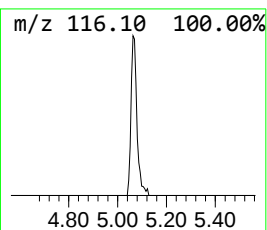
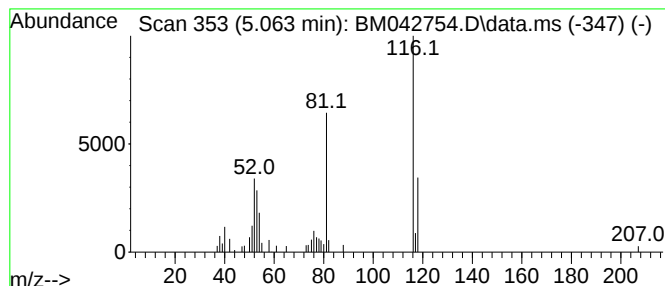
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	3.21 ng/ul	37862	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Chloro-1-methylimidazole	116	C4H5ClN2	000872-49-1	38
2			2-Furanmethanol, acetate	140	C7H8O3	000623-17-6	35
3			3-Imidazol-1-ylpropanenitrile	121	C6H7N3	023996-53-4	25
4			2-Furanmethanol, acetate	140	C7H8O3	000623-17-6	25
5			2-Furanmethanol, acetate	140	C7H8O3	000623-17-6	25



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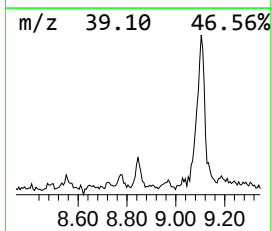
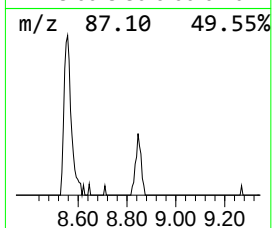
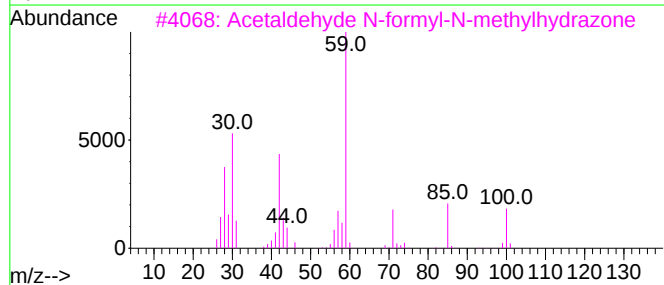
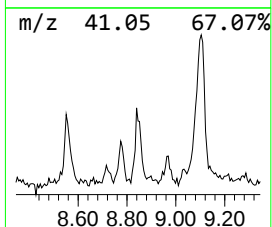
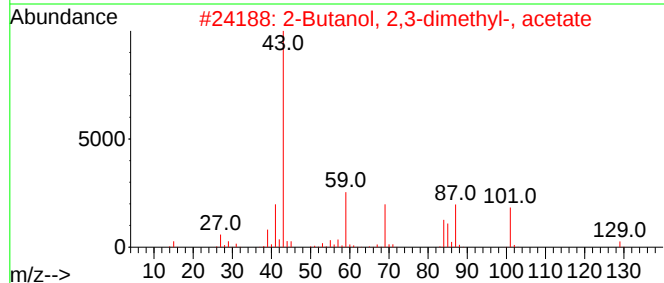
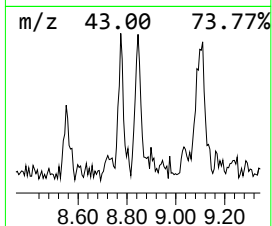
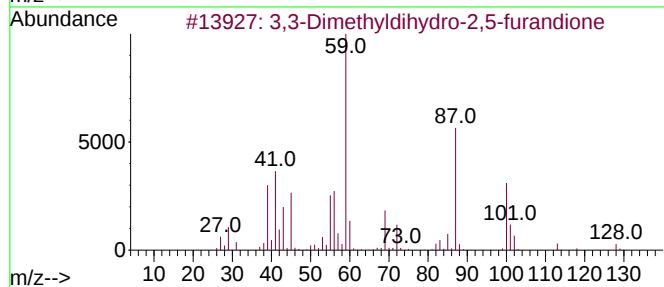
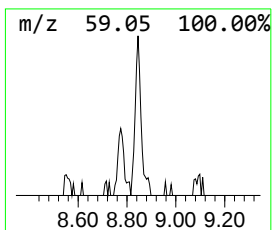
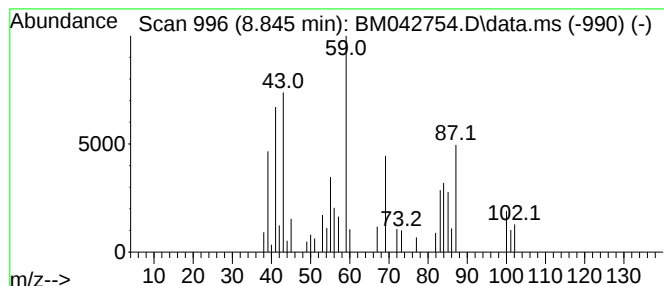
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 3,3-Dimethyldihydro-2,5-fur... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.845	2.19 ng/ul	25792	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,3-Dimethyldihydro-2,5-furandione	128	C6H8O3	017347-61-4	50
2			2-Butanol, 2,3-dimethyl-, acetate	144	C8H16O2	004806-33-1	37
3			Acetaldehyde N-formyl-N-methylhy...	100	C4H8N2O	016568-02-8	32
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	27
5			2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	25



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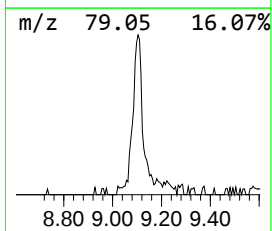
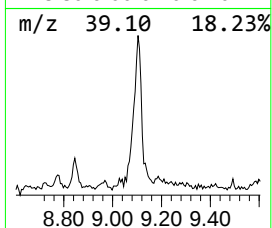
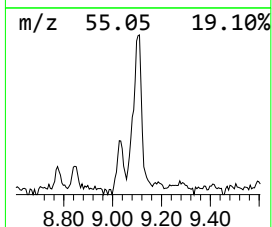
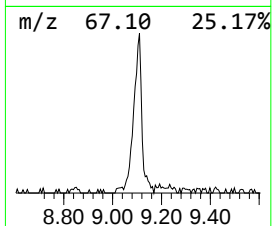
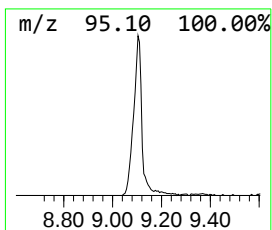
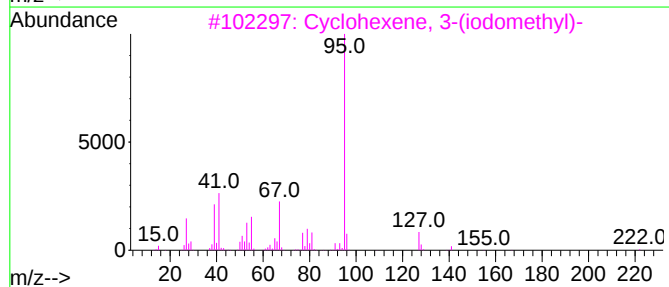
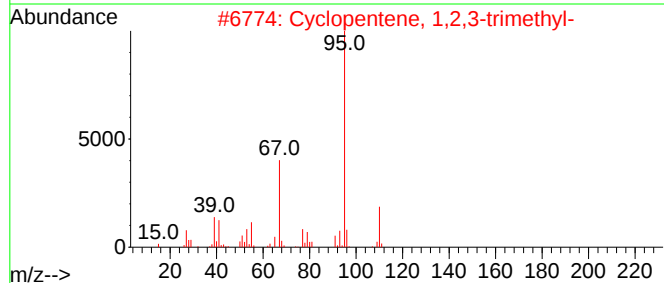
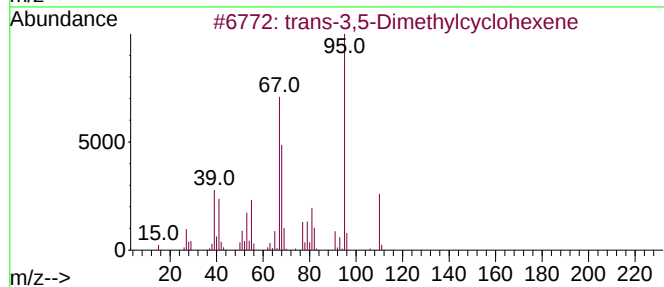
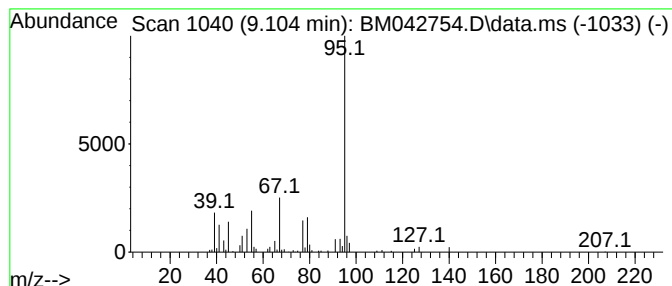
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 trans-3,5-Dimethylcyclohexene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.104	19.64 ng/ul	231622	1,4-Dichlorobenzene-d4	7.834	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-3,5-Dimethylcyclohexene	110	C8H14	056021-63-7	64
2	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	59
3	Cyclohexene, 3-(iodomethyl)-	222	C7H11I	034825-94-0	56
4	1,4-Pentadiene, 2,3,4-trimethyl-	110	C8H14	072014-90-5	50
5	1,2-Dimethyl-cyclopent-2-enecarb...	140	C8H12O2	1000190-58-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
Operator : MA/JU
Sample : 05079-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHEF1

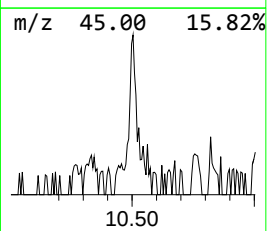
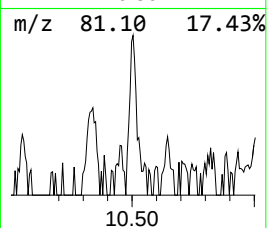
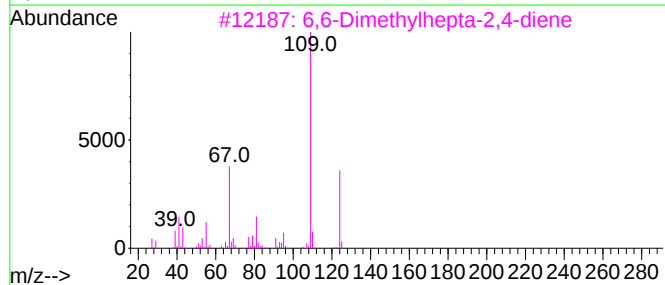
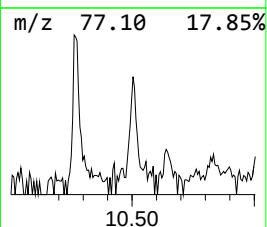
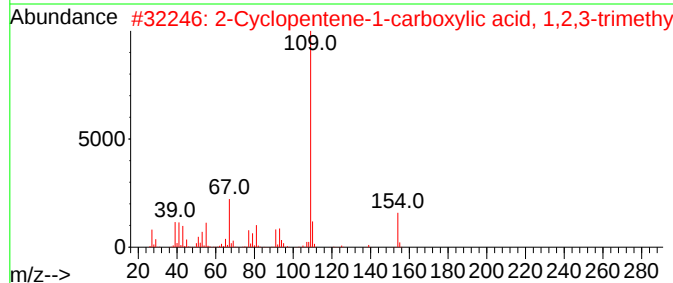
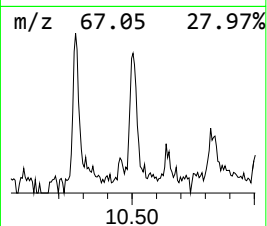
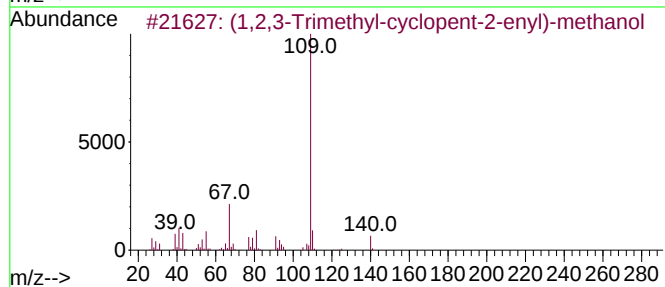
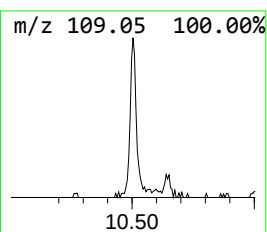
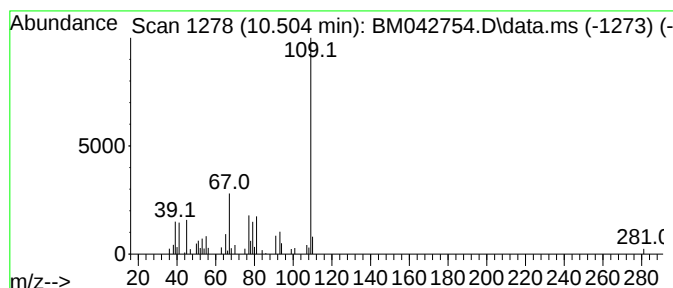
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (1,2,3-Trimethyl-cyclopent-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.504	3.62 ng/ul	60003	Naphthalene-d8	10.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1,2,3-Trimethyl-cyclopent-2-eny...	140	C9H16O	1000190-91-3	74
2			2-Cyclopentene-1-carboxylic acid...	154	C9H14O2	006894-69-5	72
3			6,6-Dimethylhepta-2,4-diene	124	C9H16	1000195-03-3	72
4			Ethanone, 1-(1,3-dimethyl-3-cycl...	152	C10H16O	051733-68-7	72
5			3,3,5,5-Tetramethylcyclopentene	124	C9H16	038667-10-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
Operator : MA/JU
Sample : 05079-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHEF1

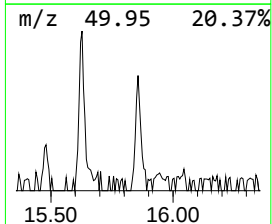
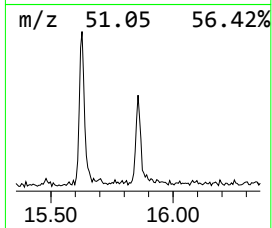
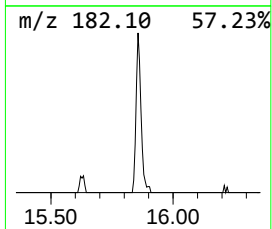
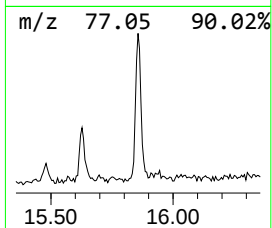
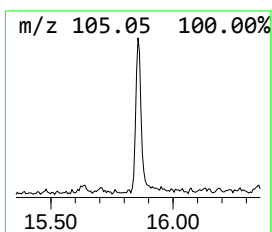
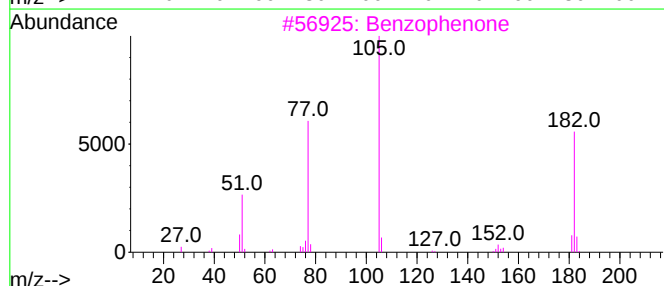
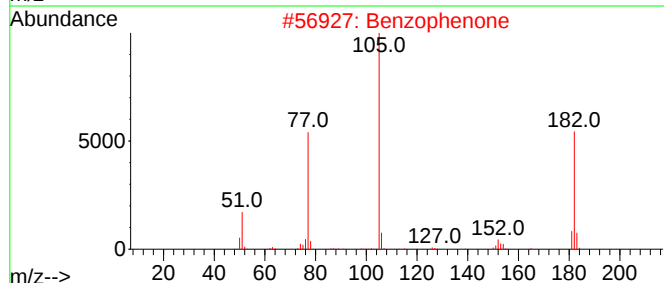
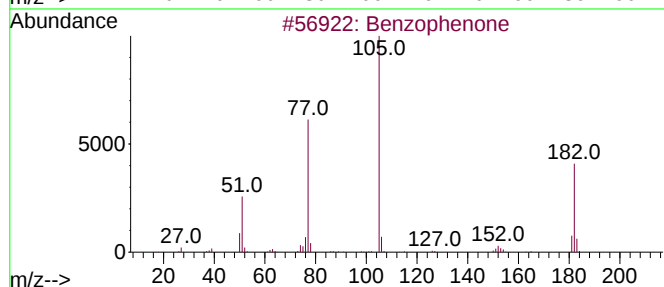
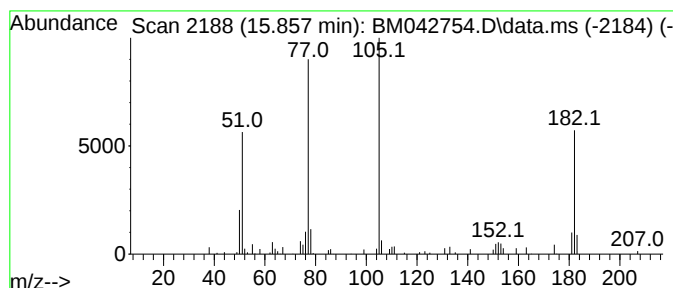
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.857	2.06 ng/ul	45739	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	91
2			Benzophenone	182	C13H10O	000119-61-9	91
3			Benzophenone	182	C13H10O	000119-61-9	91
4			Benzophenone	182	C13H10O	000119-61-9	86
5			Benzophenone	182	C13H10O	000119-61-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
Operator : MA/JU
Sample : 05079-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

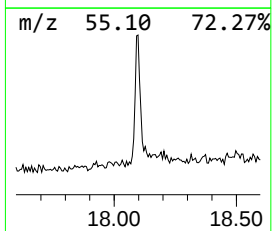
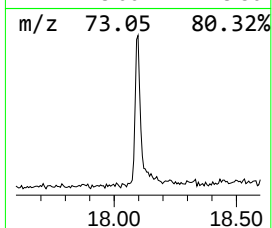
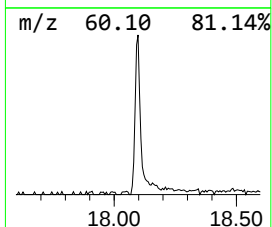
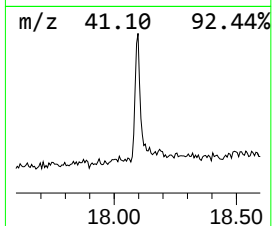
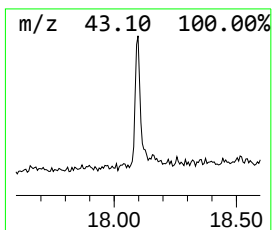
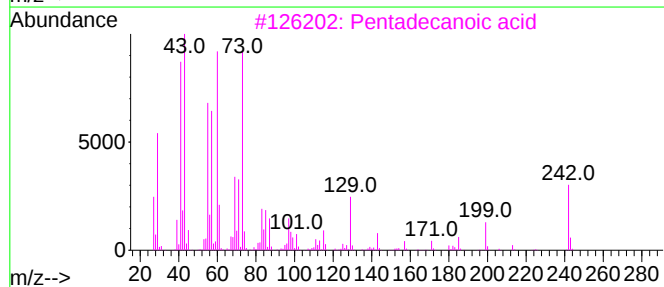
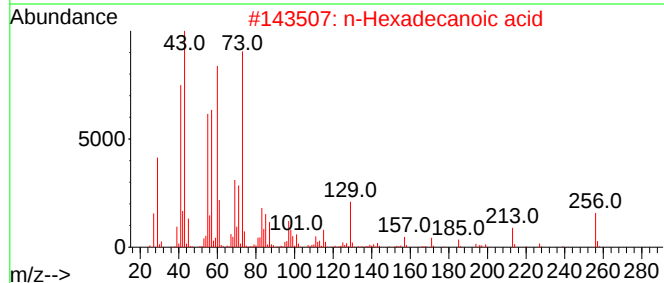
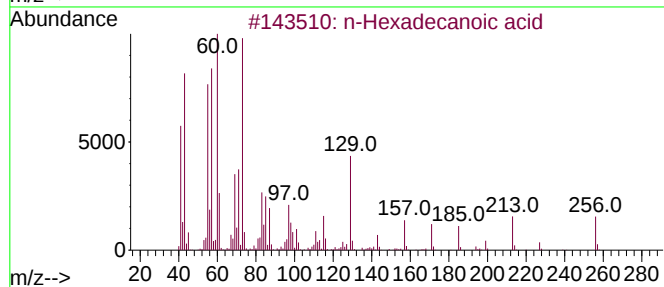
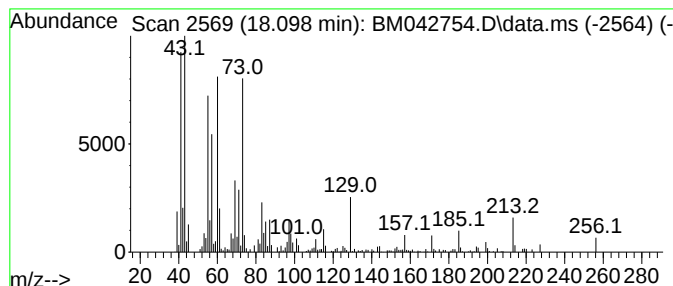
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.098	8.16 ng/ul	218275	Phenanthrene-d10	17.239	
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	87
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4	Tridecanoic acid	214	C13H26O2	000638-53-9	81
5	Tridecanoic acid	214	C13H26O2	000638-53-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
Operator : MA/JU
Sample : 05079-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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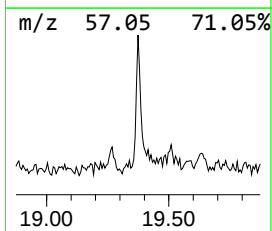
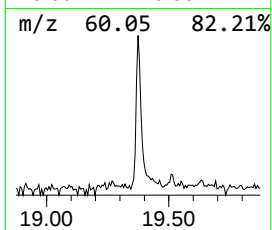
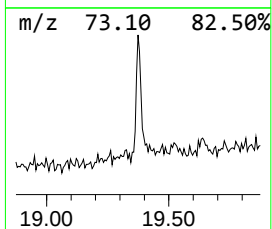
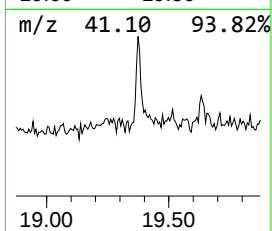
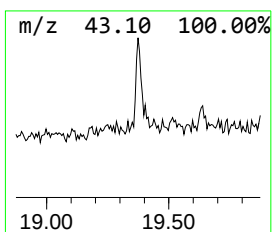
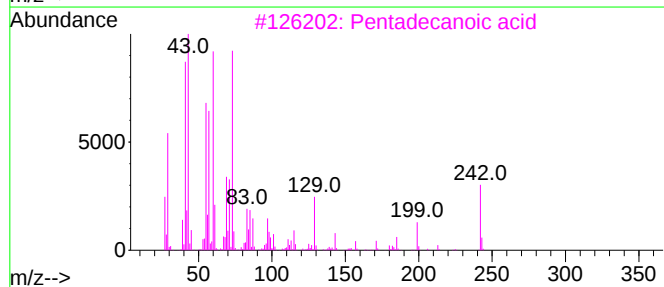
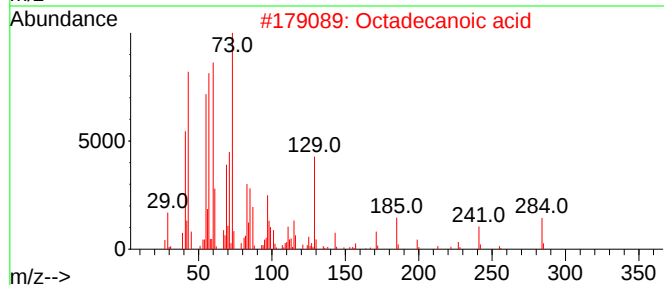
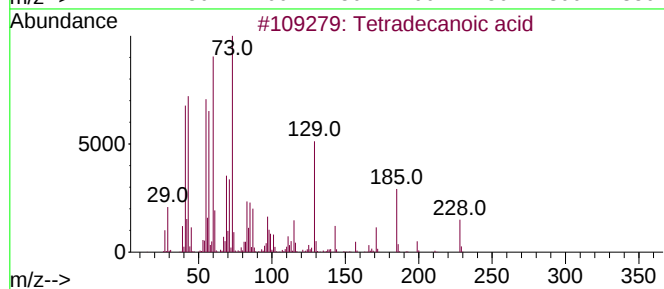
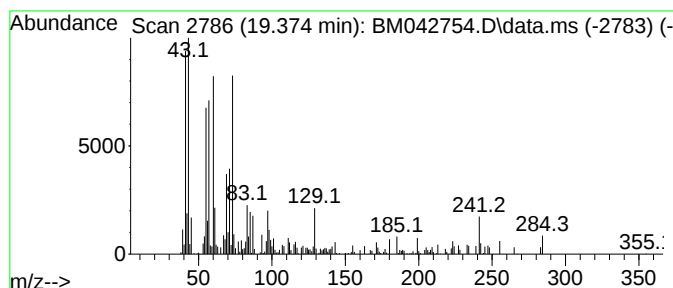
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tetradecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.374	3.37 ng/ul	95370	Chrysene-d12	21.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
2			Octadecanoic acid	284	C18H36O2	000057-11-4	83
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
Operator : MA/JU
Sample : 05079-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

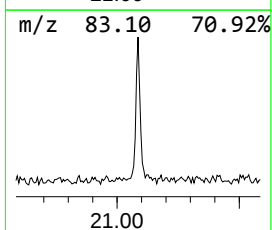
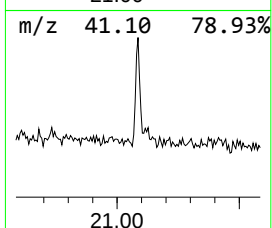
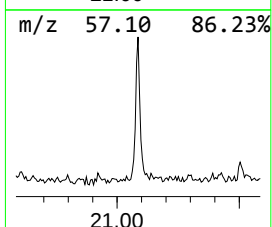
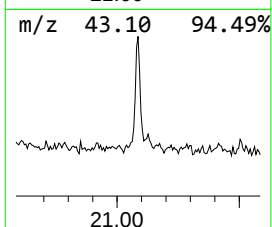
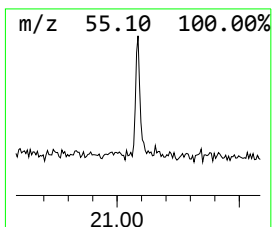
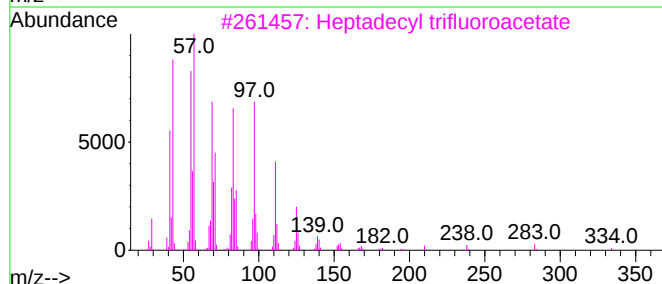
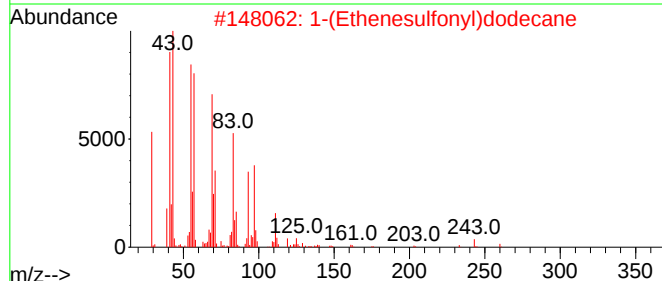
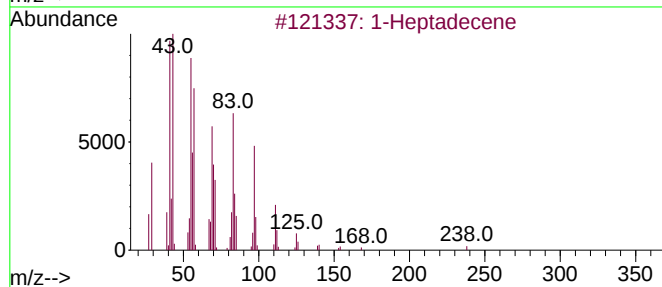
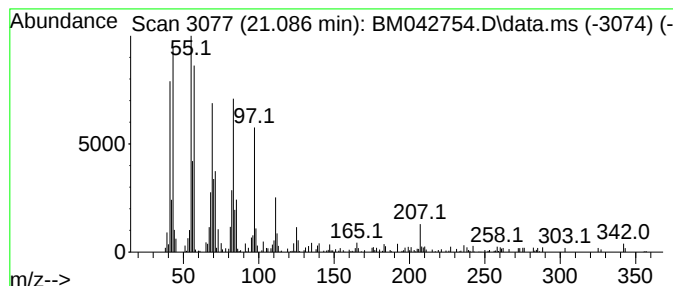
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.086	3.67 ng/ul	103963	Chrysene-d12	21.421	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptadecene	238	C17H34	006765-39-5	92
2	1-(Ethenesulfonyl)dodecane	260	C14H28O2S	030719-66-5	91
3	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	90
4	1-Heneicosanol	312	C21H44O	015594-90-8	90
5	1-Heptacosanol	396	C27H56O	002004-39-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042754.D
Acq On : 15 Nov 2023 03:59
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ALS Vial : 57 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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

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trans-3,5-Dimet...	9.104	19.6	ng/ul	231622	1	7.834	235896	20.0
(1,2,3-Trimethy...	10.504	3.6	ng/ul	60003	2	10.645	331574	20.0
Benzophenone	15.857	2.1	ng/ul	45739	3	14.486	443204	20.0
n-Hexadecanoic ...	18.098	8.2	ng/ul	218275	4	17.239	535233	20.0
Tetradecanoic acid	19.374	3.4	ng/ul	95370	5	21.421	566566	20.0
(DEL) Alkane: S...	21.086	3.7	ng/ul	103963	5	21.421	566566	20.0