

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
 Data File : BM044354.D
 Acq On : 27 Feb 2024 05:05
 Operator : MA/JU
 Sample : P1469-17
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH9R8

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-BM022324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM044354.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	7	10	20	rVB	408491	528291	14.89%	1.002%
2	3.363	44	47	53	rVB	46535	55137	1.55%	0.105%
3	3.787	115	119	127	rBV2	347570	575474	16.22%	1.091%
4	3.893	132	137	145	rVV	293128	379755	10.70%	0.720%
5	3.963	146	149	157	rVB	44735	60707	1.71%	0.115%
6	4.104	168	173	179	rVB3	32000	53359	1.50%	0.101%
7	4.328	207	211	218	rVB	43287	64285	1.81%	0.122%
8	4.481	234	237	242	rVB3	13559	17873	0.50%	0.034%
9	4.628	258	262	264	rBV3	14168	18061	0.51%	0.034%
10	4.669	264	269	275	rVV	156741	215355	6.07%	0.408%
11	4.740	275	281	288	rVV3	25353	47639	1.34%	0.090%
12	4.822	292	295	299	rVV4	16183	25778	0.73%	0.049%
13	4.863	299	302	307	rVV5	16427	25192	0.71%	0.048%
14	4.928	309	313	319	rVB7	18117	29754	0.84%	0.056%
15	5.022	325	329	340	rVB	48054	83744	2.36%	0.159%
16	5.128	343	347	356	rBV	66154	101790	2.87%	0.193%
17	5.940	477	485	491	rBV6	14334	46611	1.31%	0.088%
18	6.793	623	630	636	rBV9	9863	21185	0.60%	0.040%
19	7.098	674	682	701	rBV	630090	1118451	31.52%	2.121%
20	7.275	706	712	721	rBV	550974	877752	24.73%	1.664%
21	7.463	737	744	753	rBV	645628	1033043	29.11%	1.959%
22	7.551	754	759	764	rVB6	10361	17316	0.49%	0.033%
23	7.934	816	824	833	rBV	943148	1498389	42.22%	2.841%
24	8.651	939	946	957	rBV	355293	629887	17.75%	1.194%
25	8.769	959	966	975	rVV	198670	334052	9.41%	0.633%
26	8.892	982	987	992	rVB4	7708	14921	0.42%	0.028%
27	9.104	1017	1023	1031	rBV	579371	964116	27.17%	1.828%
28	9.210	1038	1041	1047	rVV6	6964	14874	0.42%	0.028%
29	9.269	1047	1051	1057	rVB7	10669	18764	0.53%	0.036%
30	9.492	1084	1089	1094	rBV2	16191	24786	0.70%	0.047%
31	9.828	1137	1146	1152	rBV	450677	791052	22.29%	1.500%
32	9.875	1152	1154	1159	rVB5	13997	23997	0.68%	0.045%
33	9.975	1166	1171	1180	rBV	44133	92949	2.62%	0.176%
34	10.192	1203	1208	1212	rBV2	23716	36635	1.03%	0.069%
35	10.363	1230	1237	1247	rVV	685703	1198816	33.78%	2.273%
36	10.457	1247	1253	1264	rVV	72311	156722	4.42%	0.297%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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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-BM022324.MA.M
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37	10.739	1293	1301	1312	rBV	1313385	2179212	61.41%	4.132%
38	10.892	1321	1327	1330	rBV	123118	223248	6.29%	0.423%
39	10.933	1330	1334	1342	rVB	219061	381630	10.75%	0.724%
40	11.139	1364	1369	1376	rBV10	8557	20845	0.59%	0.040%
41	11.298	1390	1396	1403	rBV10	19818	45438	1.28%	0.086%
42	11.545	1432	1438	1446	rVV6	24996	59316	1.67%	0.112%
43	12.122	1530	1536	1540	rBV	11947	19766	0.56%	0.037%
44	12.327	1564	1571	1578	rVB3	16037	35367	1.00%	0.067%
45	13.410	1748	1755	1760	rVB3	13977	24757	0.70%	0.047%
46	13.480	1760	1767	1775	rBV	52538	99373	2.80%	0.188%
47	13.727	1802	1809	1813	rBV10	9365	23782	0.67%	0.045%
48	13.827	1820	1826	1830	rBV6	43367	78482	2.21%	0.149%
49	13.863	1830	1832	1835	rVV3	22343	26971	0.76%	0.051%
50	13.998	1846	1855	1859	rBV	1293711	1884444	53.10%	3.573%
51	14.033	1859	1861	1868	rVV	218663	263832	7.43%	0.500%
52	14.104	1868	1873	1880	rVB	779363	1123432	31.66%	2.130%
53	14.169	1880	1884	1890	rBV7	12106	16359	0.46%	0.031%
54	14.269	1890	1901	1912	rBV	1477269	2291989	64.58%	4.345%
55	14.363	1912	1917	1922	rVB4	51714	84663	2.39%	0.161%
56	14.439	1923	1930	1934	rBV4	19059	28823	0.81%	0.055%
57	14.574	1947	1953	1962	rBV	1810687	2705296	76.23%	5.129%
58	14.786	1982	1989	2000	rBV	352389	715187	20.15%	1.356%
59	14.892	2002	2007	2013	rVV3	138950	296618	8.36%	0.562%
60	14.945	2013	2016	2022	rVV2	67971	126511	3.56%	0.240%
61	15.004	2022	2026	2031	rVV6	64391	139014	3.92%	0.264%
62	15.057	2033	2035	2040	rVB6	26048	31786	0.90%	0.060%
63	15.298	2072	2076	2083	rVV2	23245	43927	1.24%	0.083%
64	15.380	2085	2090	2095	rVV2	17920	30343	0.85%	0.058%
65	15.568	2112	2122	2131	rBV	2048183	2938879	82.81%	5.572%
66	15.692	2137	2143	2152	rVV	505624	710192	20.01%	1.346%
67	15.798	2156	2161	2166	rBV6	15000	33043	0.93%	0.063%
68	16.298	2242	2246	2251	rVB	62828	84138	2.37%	0.160%
69	16.368	2253	2258	2266	rVV2	46491	91203	2.57%	0.173%
70	16.439	2266	2270	2277	rVB10	16402	31633	0.89%	0.060%
71	16.545	2281	2288	2300	rBV	277719	427649	12.05%	0.811%
72	16.733	2316	2320	2327	rBV6	32563	48095	1.36%	0.091%
73	16.874	2340	2344	2350	rVB2	26454	37884	1.07%	0.072%
74	17.051	2370	2374	2377	rBV2	40972	55276	1.56%	0.105%
75	17.098	2377	2382	2393	rVB	286710	382958	10.79%	0.726%
76	17.204	2396	2400	2402	rBV	102820	125635	3.54%	0.238%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	17.233	2402	2405	2412	rVV2	108707	159055	4.48%	0.302%
78	17.327	2412	2421	2432	rVV	2151084	3162969	89.13%	5.997%
79	17.427	2432	2438	2448	rVV	1953278	2834408	79.87%	5.374%
80	17.509	2448	2452	2459	rVB8	22451	46836	1.32%	0.089%
81	17.721	2485	2488	2495	rBV9	13346	23816	0.67%	0.045%
82	17.839	2504	2508	2513	rVB5	22743	30720	0.87%	0.058%
83	18.009	2535	2537	2542	rVB6	14528	18508	0.52%	0.035%
84	18.115	2549	2555	2561	rBV3	94265	170657	4.81%	0.324%
85	18.180	2561	2566	2571	rBV2	100368	148990	4.20%	0.282%
86	18.239	2571	2576	2587	rBV	973833	1456675	41.05%	2.762%
87	18.533	2622	2626	2630	rBV4	57659	100715	2.84%	0.191%
88	18.674	2646	2650	2654	rBV2	68734	100718	2.84%	0.191%
89	18.809	2669	2673	2683	rVB	83398	153361	4.32%	0.291%
90	19.039	2709	2712	2716	rBV	40495	49029	1.38%	0.093%
91	19.080	2716	2719	2727	rVB7	43746	63848	1.80%	0.121%
92	19.398	2771	2773	2776	rBV3	43756	48749	1.37%	0.092%
93	19.521	2790	2794	2803	rBV	281843	400628	11.29%	0.760%
94	19.721	2823	2828	2836	rBV	2544524	3548895	100.00%	6.729%
95	20.674	2987	2990	2993	rVB	83380	96882	2.73%	0.184%
96	21.256	3085	3089	3097	rBV	511977	708867	19.97%	1.344%
97	21.527	3130	3135	3140	rBV	2271880	3077741	86.72%	5.835%
98	23.168	3409	3414	3421	rVB	828353	1272478	35.86%	2.413%
99	23.786	3513	3519	3528	rBV	1504977	2913257	82.09%	5.523%
100	23.944	3540	3546	3553	rVB	1471204	2992686	84.33%	5.674%

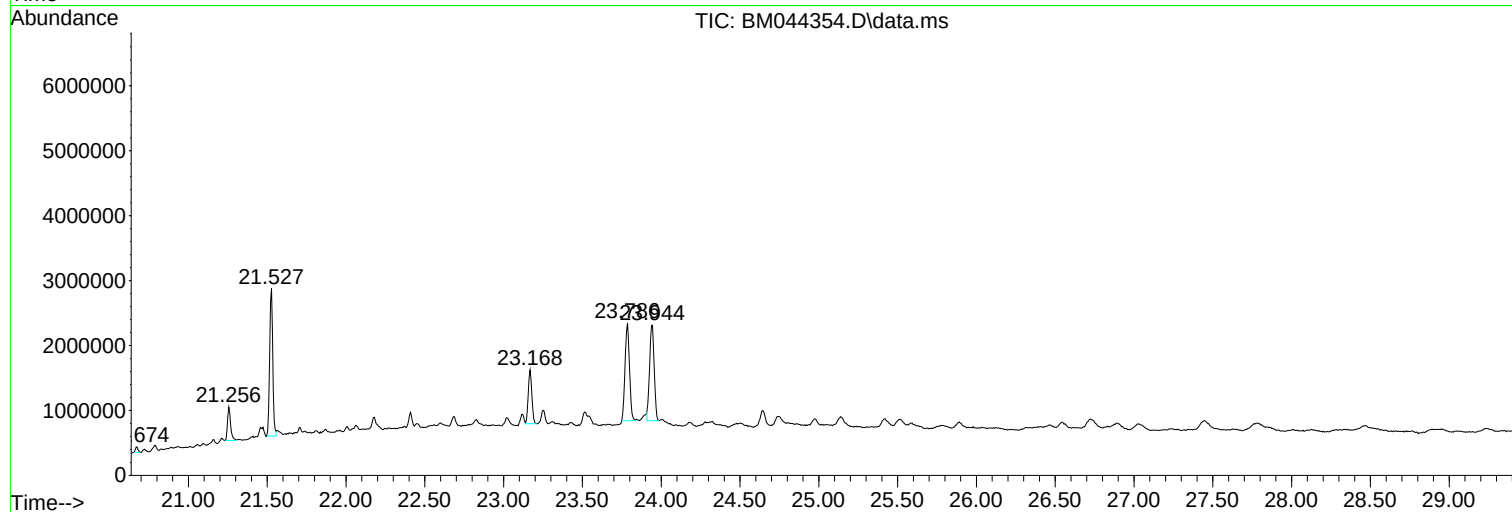
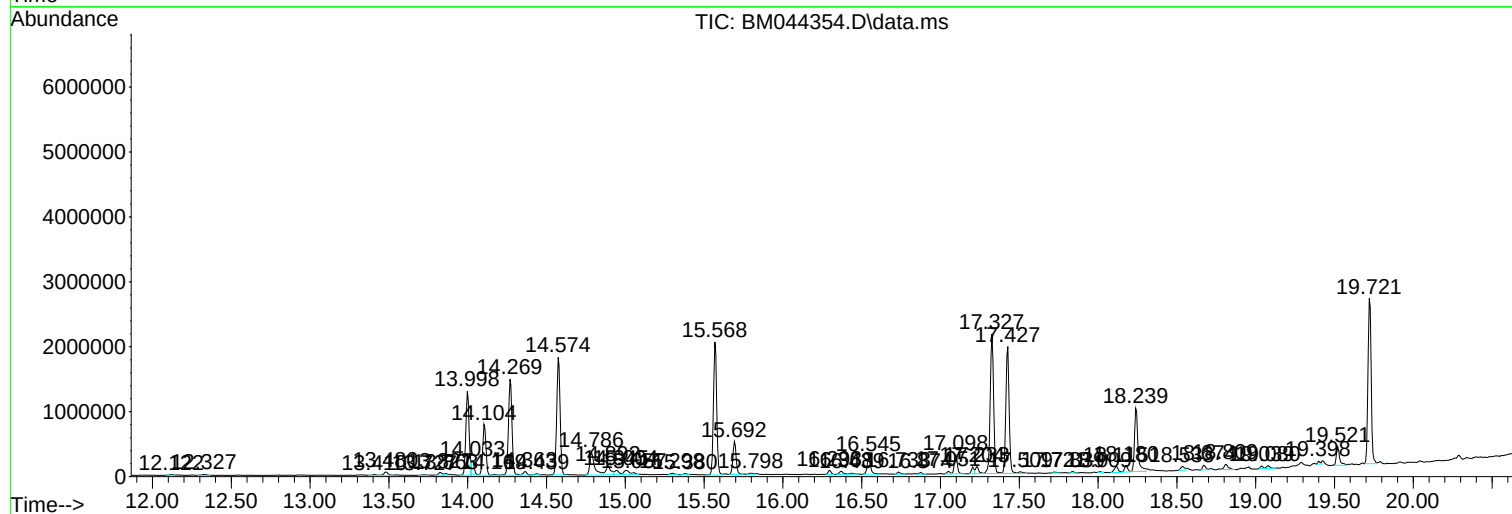
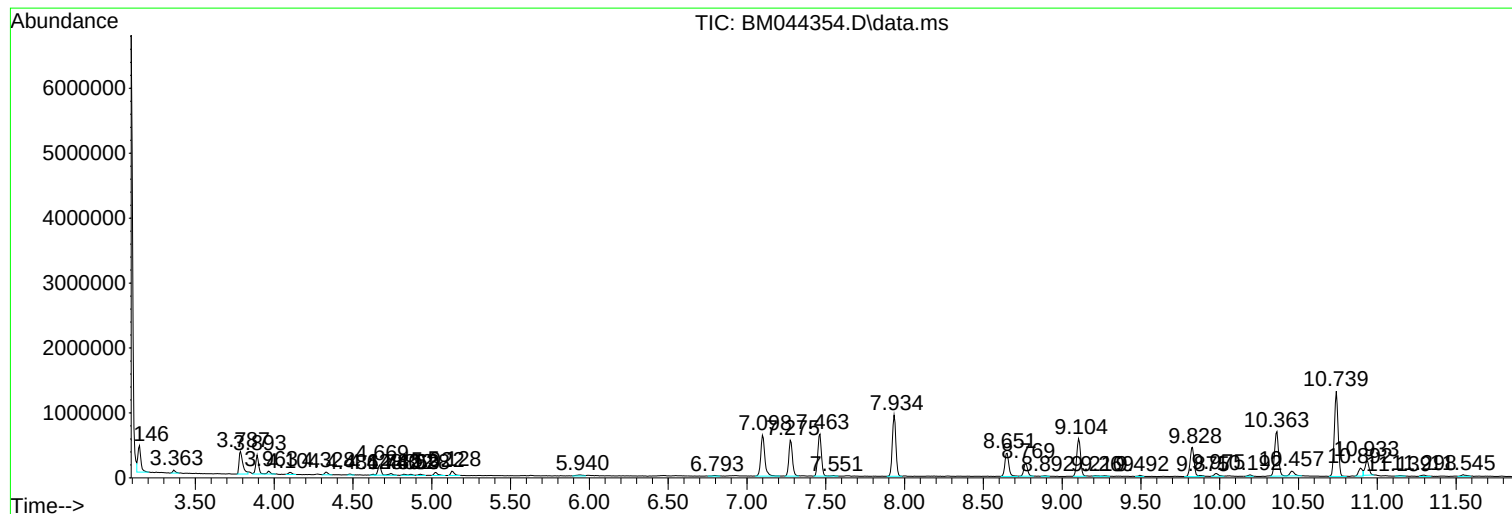
Sum of corrected areas: 52743996

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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



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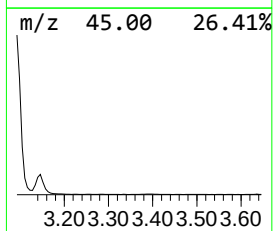
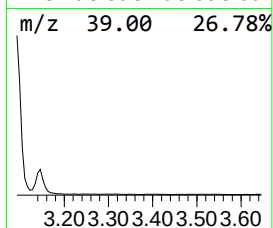
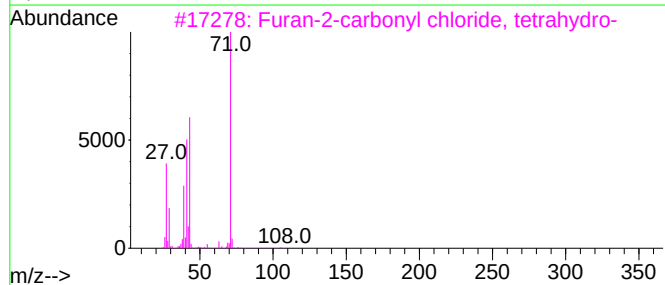
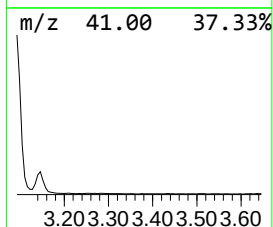
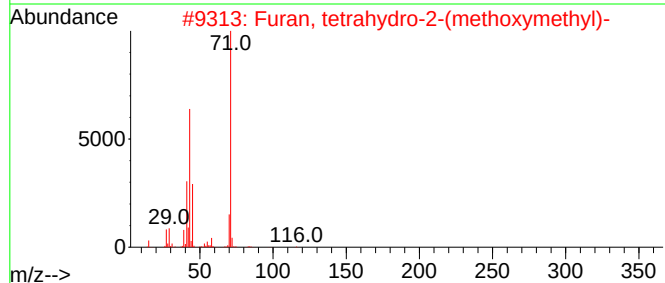
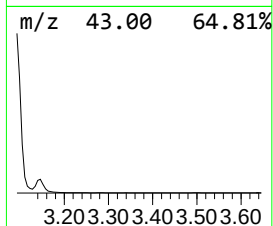
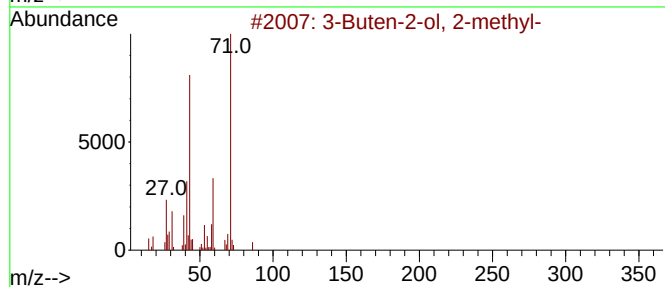
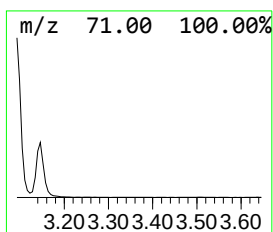
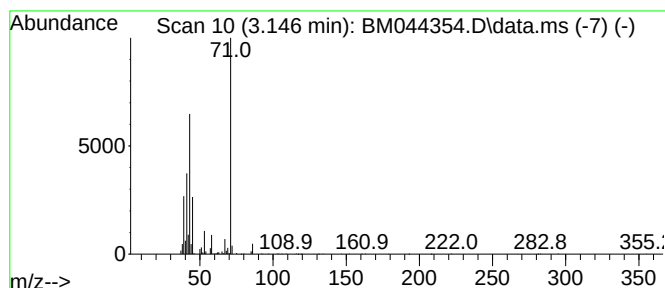
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Buten-2-ol, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.146	7.05 ng/ul	528291	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	80
2			Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	64
3			Furan-2-carbonyl chloride, tetra...	134	C5H7ClO2	052449-98-6	47
4			3-Penten-2-ol	86	C5H10O	001569-50-2	46
5			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	45



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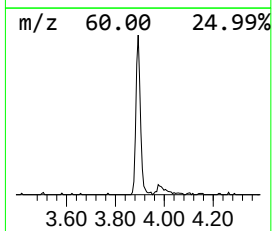
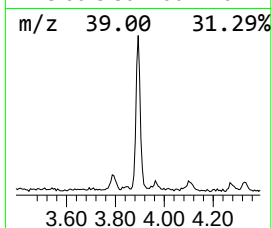
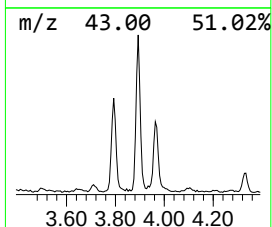
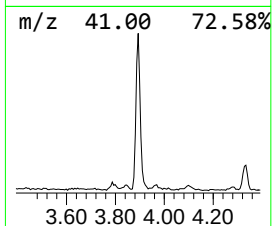
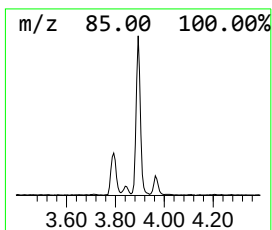
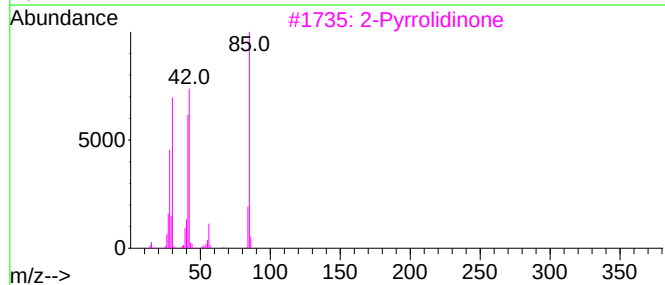
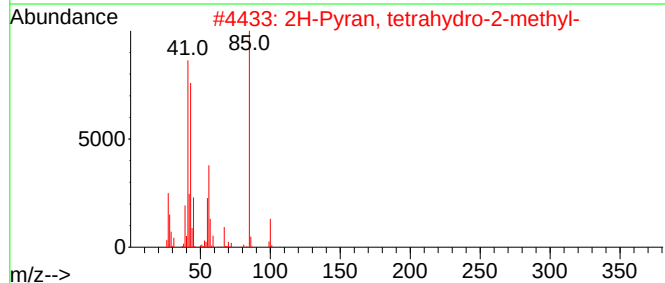
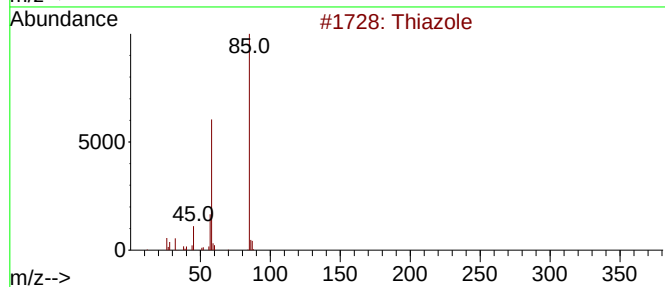
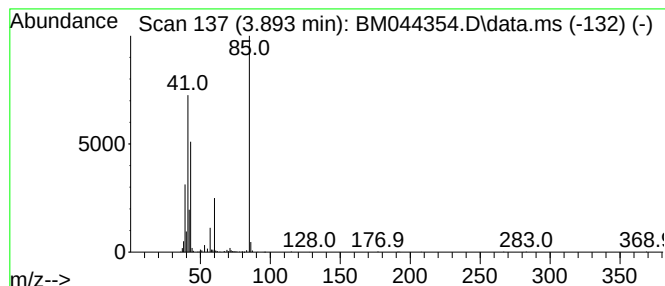
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.893	5.07 ng/ul	379755	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiazole	85	C3H3NS	000288-47-1	38
2			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3			2-Pyrrolidinone	85	C4H7NO	000616-45-5	38
4			Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	36
5			2-Pyrrolidinone	85	C4H7NO	000616-45-5	28



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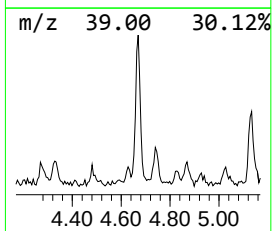
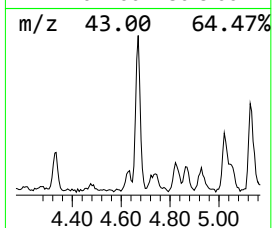
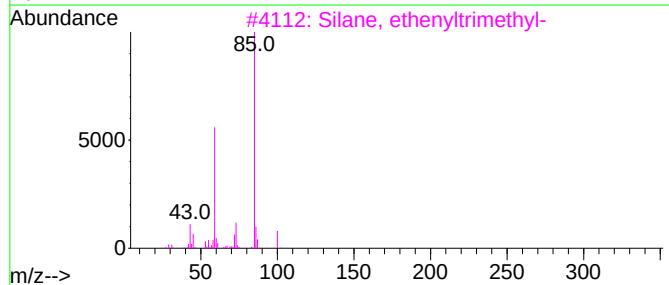
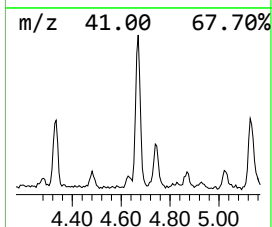
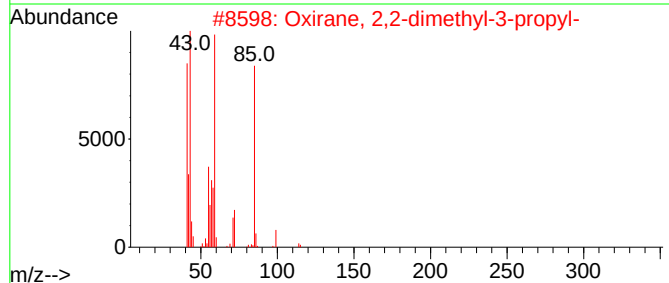
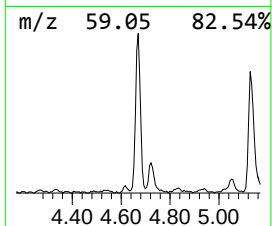
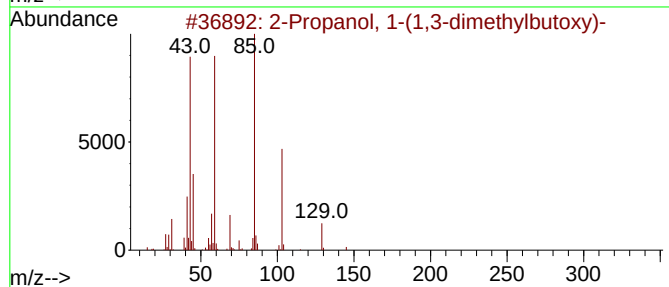
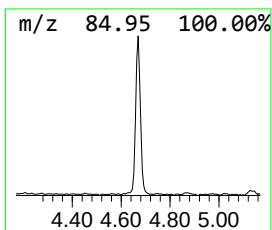
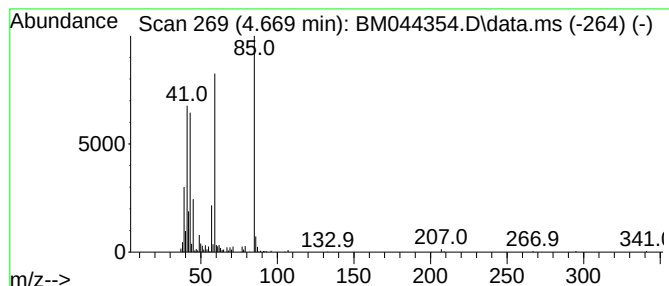
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-(1,3-dimethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.669	2.87 ng/ul	215355	1,4-Dichlorobenzene-d4	7.934

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propanol, 1-(1,3-dimethylbutoxy)-	160	C9H20O2	054340-89-5	72
2		Oxirane, 2,2-dimethyl-3-propyl-	114	C7H14O	017612-35-0	64
3		Silane, ethenyltrimethyl-	100	C5H12Si	000754-05-2	38
4		N-Chloroacetyl-dl-erythro-0-meth...	209	C7H12ClNO4	1010214-48-1	38
5		Silane, ethenyltrimethyl-	100	C5H12Si	000754-05-2	38



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Data File : BM044354.D
Acq On : 27 Feb 2024 05:05
Operator : MA/JU
Sample : P1469-17
Misc :
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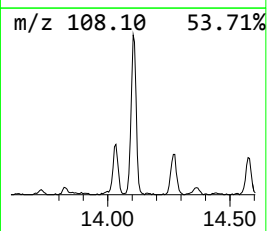
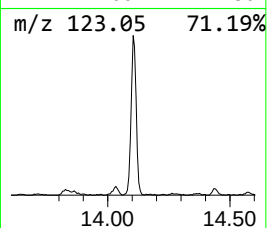
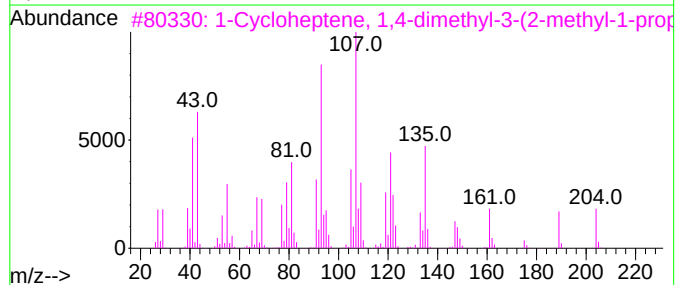
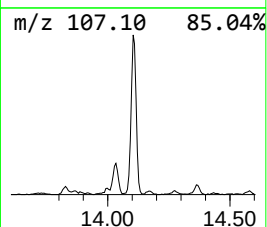
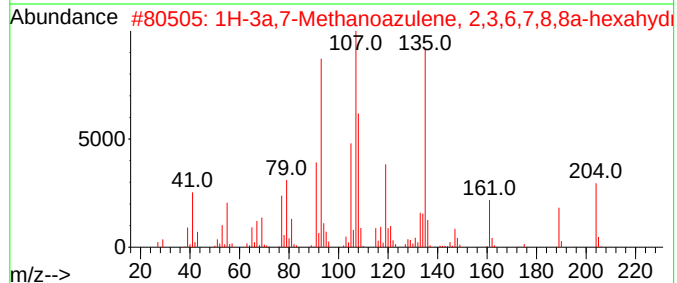
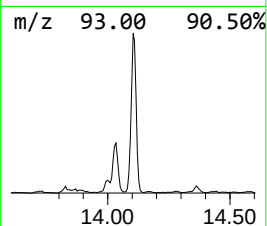
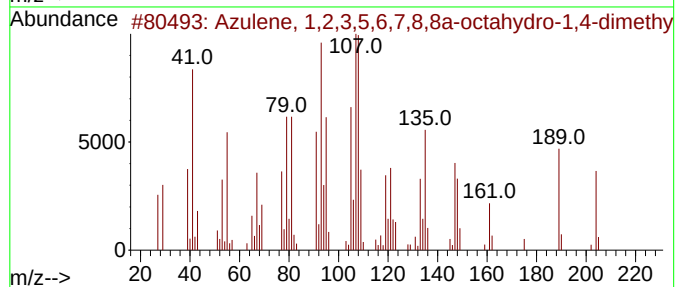
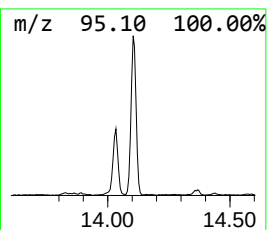
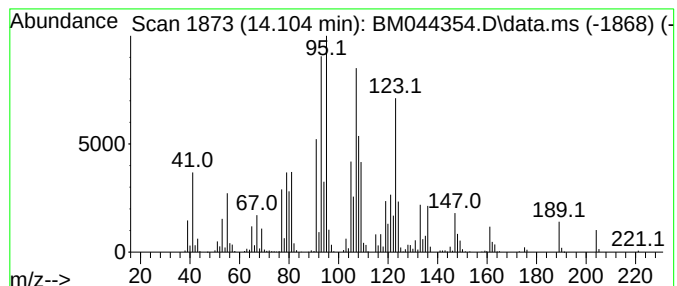
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Azulene, 1,2,3,5,6,7,8,8a-o... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.104	8.31 ng/ul	1123430	Acenaphthene-d10	14.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	58
2			1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	50
3			1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	50
4			Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	43
5			(1R,3aS,5aS,8aR)-1,3a,5a-Trimeth...	204	C15H24	071596-72-0	42



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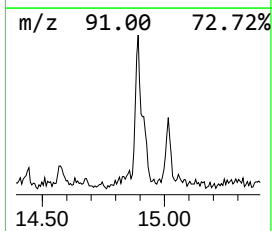
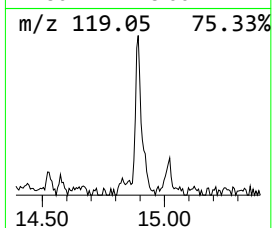
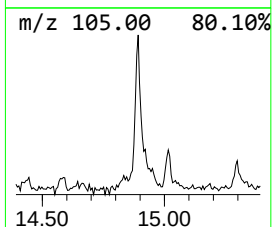
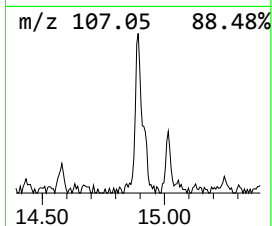
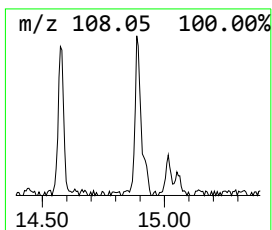
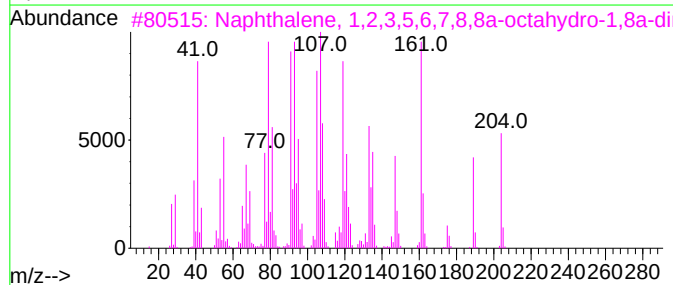
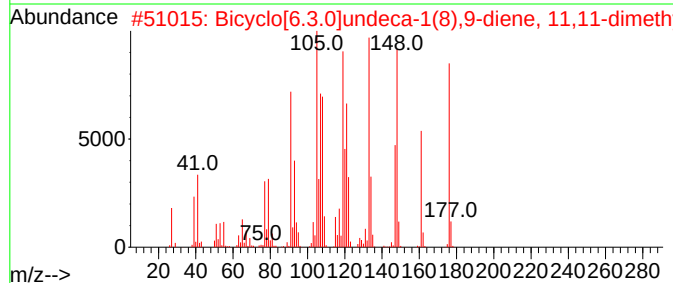
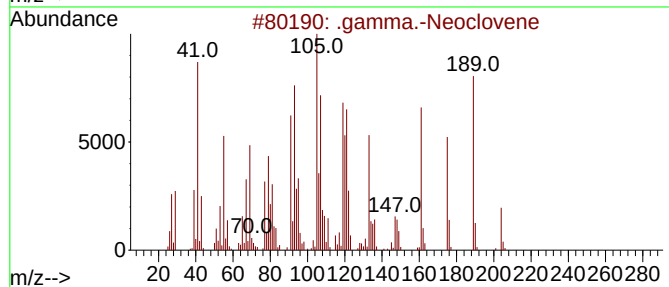
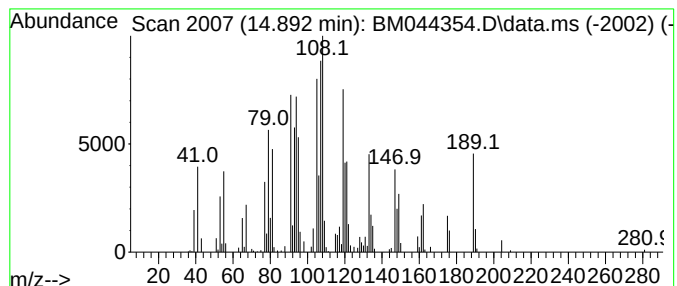
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 .gamma.-Neoclovene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.892	2.19 ng/ul	296618	Acenaphthene-d10	14.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Neoclovene	204	C15H24	1000156-11-7	64
2			Bicyclo[6.3.0]undeca-1(8),9-dien...	176	C13H20	1000163-44-3	49
3			Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	49
4			1,2,4-Methenoazulene, decahydro-...	204	C15H24	001137-12-8	45
5			Aciphyllene	204	C15H24	087745-31-1	38



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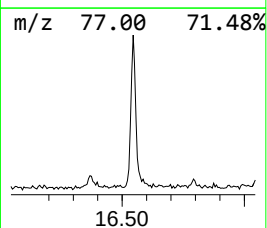
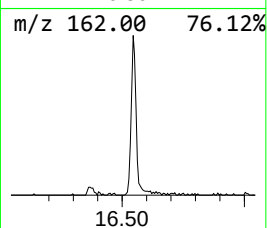
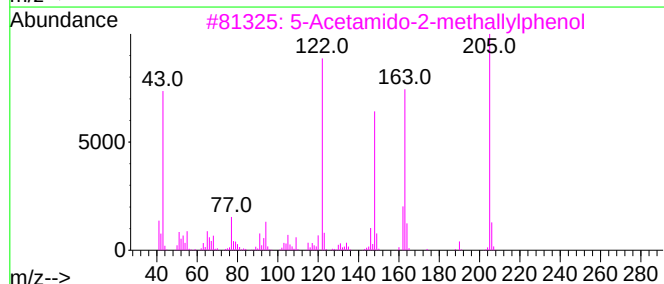
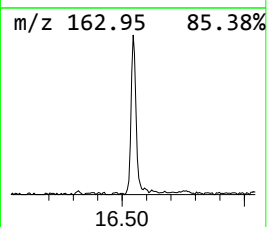
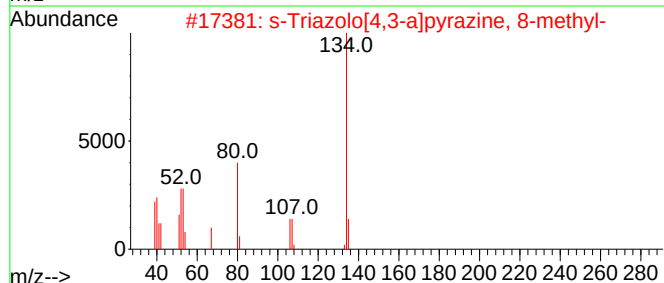
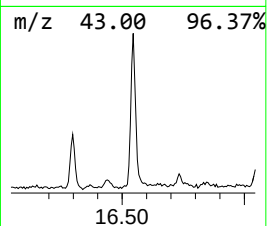
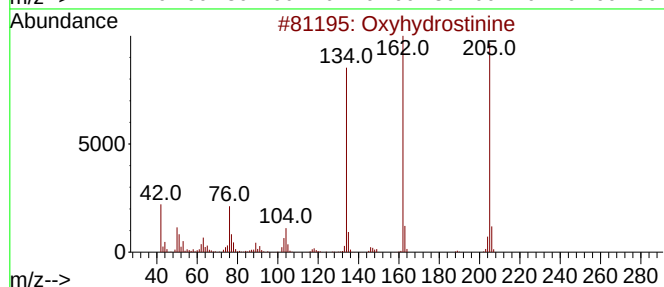
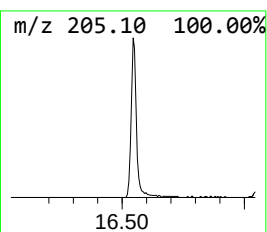
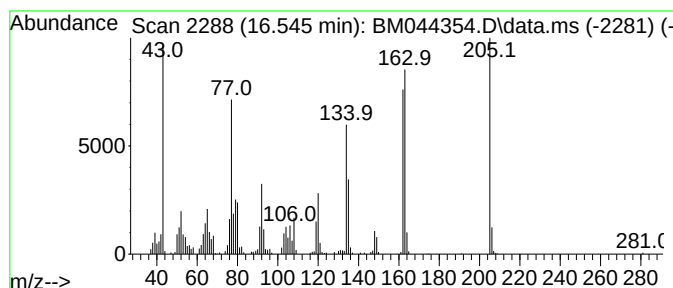
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Oxyhydrostinine Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.545	2.70 ng/ul	427649	Phenanthrene-d10	17.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxyhydrostinine	205	C11H11NO3	000552-29-4	62
2			s-Triazolo[4,3-a]pyrazine, 8-met...	134	C6H6N4	023126-45-6	35
3			5-Acetamido-2-methallylphenol	205	C12H15NO2	088913-21-7	30
4			2H-1-Benzopyran-2-one, 3,4-dihyd...	162	C10H10O2	000092-47-7	30
5			Piperidine, 1-(4-methyl-4-penten...	163	C11H17N	1000351-45-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
Data File : BM044354.D
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Misc :
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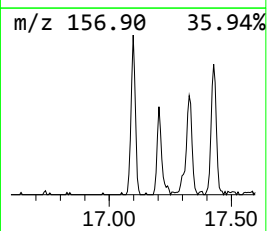
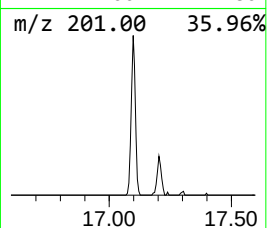
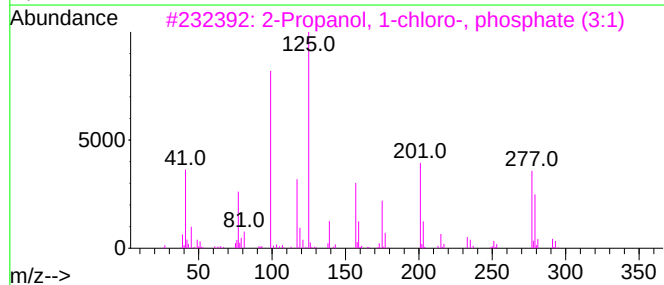
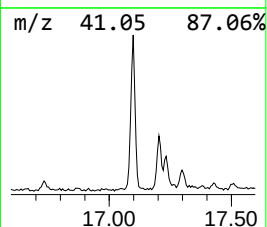
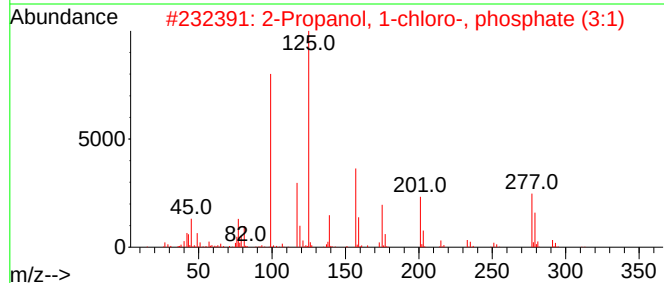
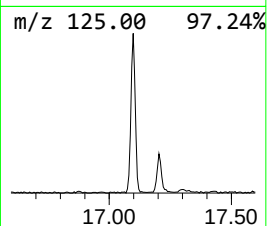
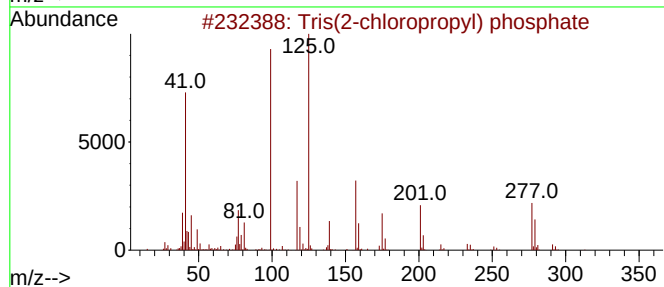
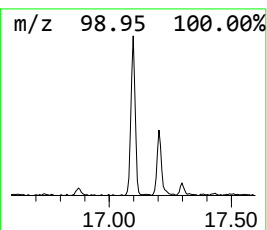
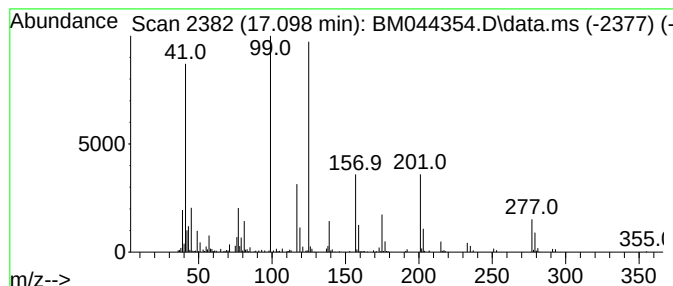
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tris(2-chloropropyl) phosphate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.098	2.42 ng/ul	382958	Phenanthrene-d10	17.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tris(2-chloropropyl) phosphate	326	C9H18Cl3O4P	006145-73-9	91
2			2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	83
3			2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	81
4			Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	43
5			Sarin	140	C4H10F02P	000107-44-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
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Misc :
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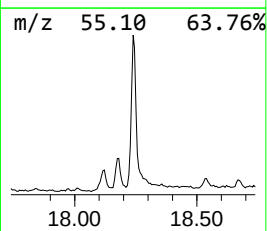
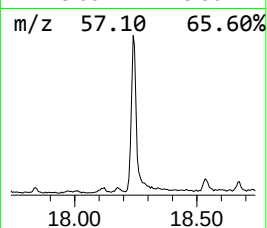
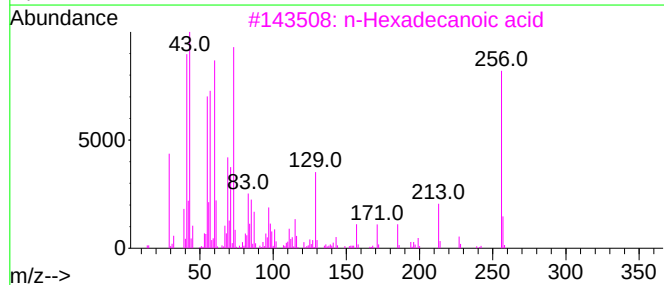
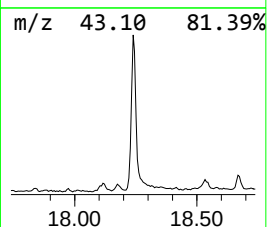
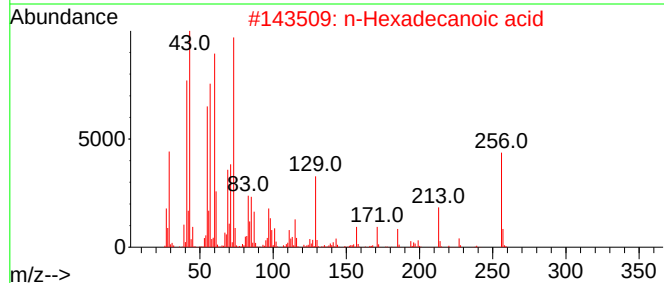
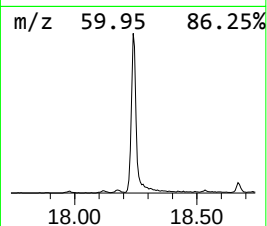
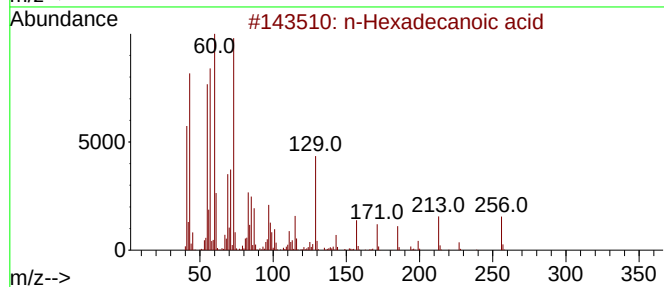
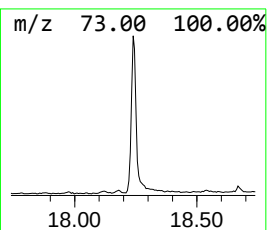
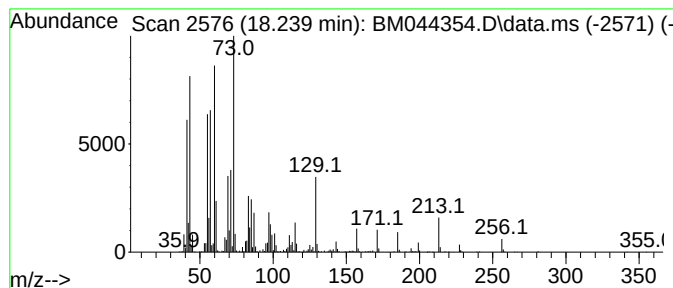
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	9.21 ng/ul	1456680	Phenanthrene-d10	17.327

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
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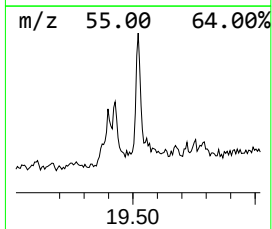
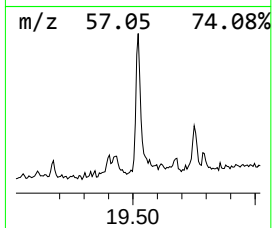
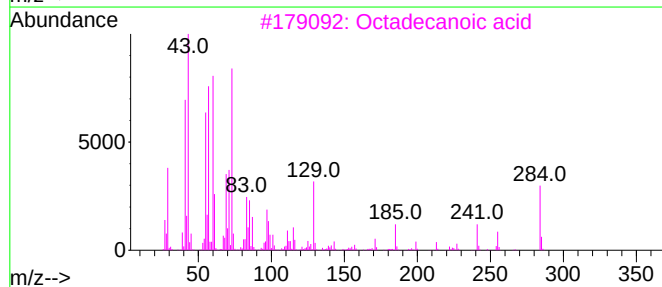
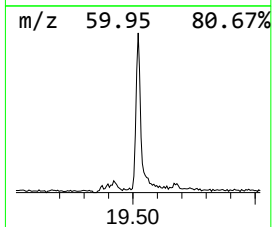
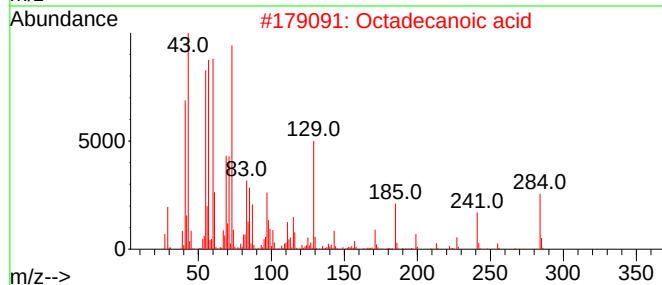
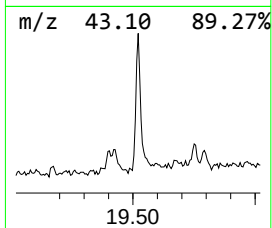
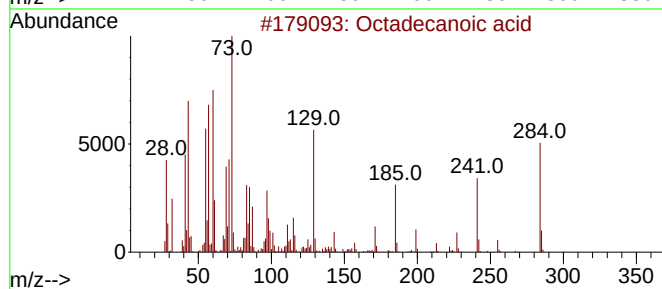
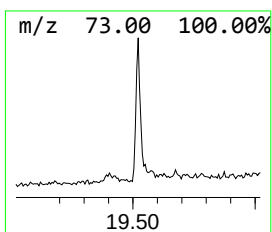
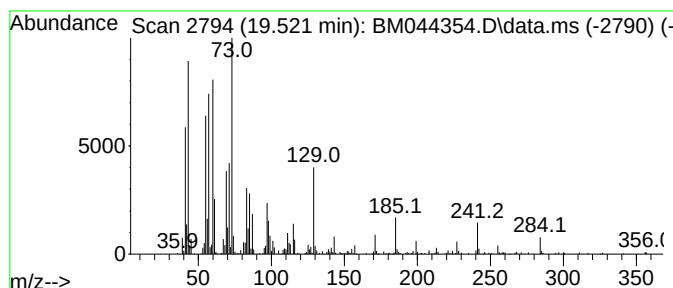
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.521	2.60 ng/ul	400628	Chrysene-d12	21.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
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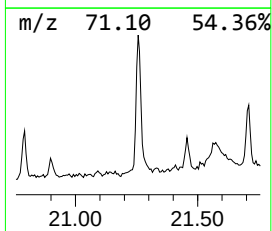
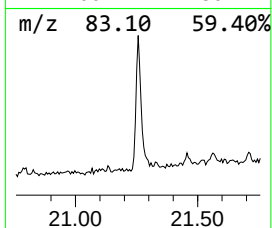
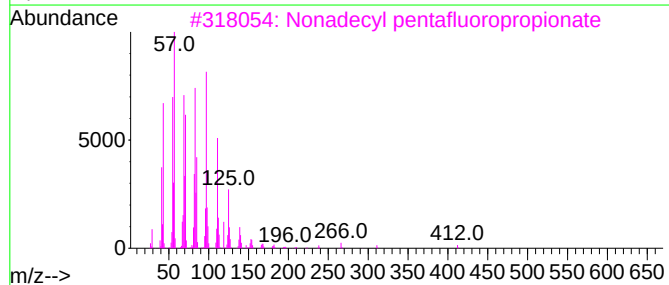
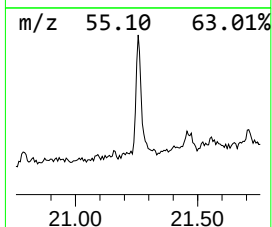
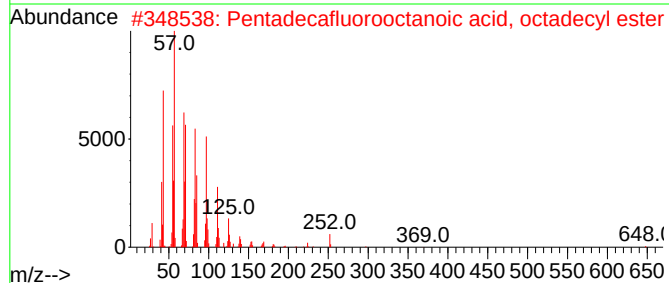
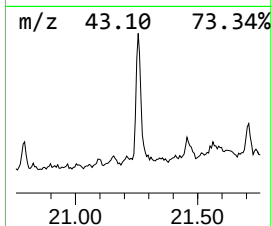
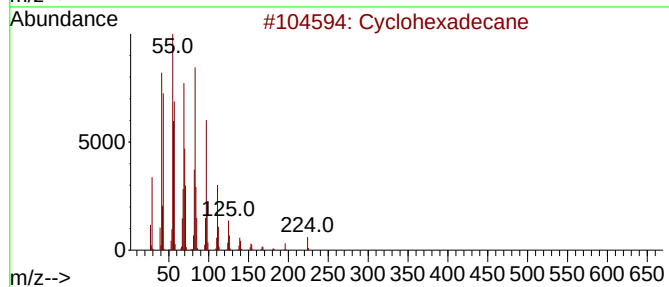
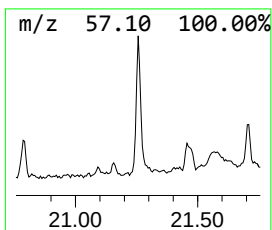
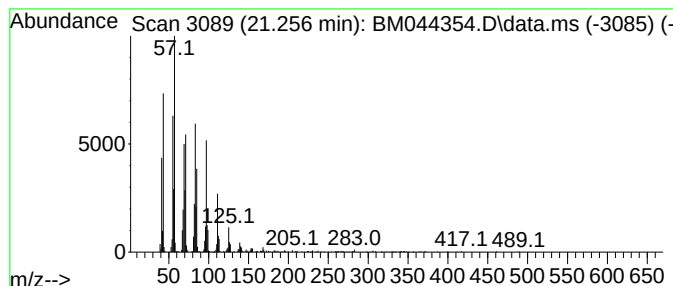
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic21.256 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.256	4.61 ng/ul	708867	Chrysene-d12	21.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	93
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
4			Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	91
5			Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
Data File : BM044354.D
Acq On : 27 Feb 2024 05:05
Operator : MA/JU
Sample : P1469-17
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH9R8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.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
3-Buten-2-ol, 2...	3.146	7.0	ng/ul	528291	1	7.934	1498390	20.0
unknown-01	3.893	5.1	ng/ul	379755	1	7.934	1498390	20.0
2-Propanol, 1-(...	4.669	2.9	ng/ul	215355	1	7.934	1498390	20.0
Azulene, 1,2,3,...	14.104	8.3	ng/ul	1123430	3	14.574	2705300	20.0
.gamma.-Neoclovene	14.892	2.2	ng/ul	296618	3	14.574	2705300	20.0
Oxyhydrostinine	16.545	2.7	ng/ul	427649	4	17.327	3162970	20.0
Tris(2-chloropr...	17.098	2.4	ng/ul	382958	4	17.327	3162970	20.0
n-Hexadecanoic ...	18.239	9.2	ng/ul	1456680	4	17.327	3162970	20.0
Octadecanoic acid	19.521	2.6	ng/ul	400628	5	21.527	3077740	20.0
(DEL) Alkane: C...	21.256	4.6	ng/ul	708867	5	21.527	3077740	20.0
unknown-02	23.168	8.5	ng/ul	1272480	6	23.944	2992690	20.0