

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033979.D
 Acq On : 02 Feb 2022 18:26
 Operator : CG/JU
 Sample : N1355-02
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COVE3

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-BM013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM033979.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.313	5	13	23	rVB	973822	1243379	4.18%	0.858%
2	3.649	65	70	87	rVV	2004241	2802052	9.41%	1.933%
3	3.878	104	109	117	rBV	479063	651273	2.19%	0.449%
4	4.060	117	140	143	rBV2	314325	1357336	4.56%	0.936%
5	4.396	194	197	199	rBV	210361	270703	0.91%	0.187%
6	4.472	206	210	215	rBV	950327	1593052	5.35%	1.099%
7	4.890	275	281	285	rVV	195701	316840	1.06%	0.219%
8	4.925	285	287	295	rVV2	80657	162167	0.54%	0.112%
9	5.049	295	308	311	rVV	9461377	21053964	70.73%	14.522%
10	5.084	311	314	322	rVB	195041	279824	0.94%	0.193%
11	5.407	363	369	377	rVV2	460158	742136	2.49%	0.512%
12	5.543	387	392	406	rVB	668768	1293654	4.35%	0.892%
13	5.754	423	428	437	rVV	309622	454530	1.53%	0.314%
14	5.925	451	457	464	rVB2	647890	1057161	3.55%	0.729%
15	6.001	464	470	482	rBV	1990806	3003612	10.09%	2.072%
16	6.190	494	502	508	rBV2	70565	145094	0.49%	0.100%
17	6.413	534	540	541	rBV3	90338	140796	0.47%	0.097%
18	6.437	541	544	548	rVV2	132790	245043	0.82%	0.169%
19	6.478	548	551	558	rVB3	143896	238971	0.80%	0.165%
20	6.560	560	565	569	rBV4	96091	183947	0.62%	0.127%
21	6.884	615	620	622	rBV2	81815	132077	0.44%	0.091%
22	6.919	622	626	630	rVB2	172542	257114	0.86%	0.177%
23	6.978	630	636	638	rBV	258233	425255	1.43%	0.293%
24	7.013	638	642	646	rVB2	304326	470732	1.58%	0.325%
25	7.066	646	651	652	rBV2	258230	333422	1.12%	0.230%
26	7.090	652	655	661	rVB2	651036	1030550	3.46%	0.711%
27	7.143	661	664	669	rVB	149649	225862	0.76%	0.156%
28	7.248	673	682	688	rBV	456012	782479	2.63%	0.540%
29	7.313	688	693	696	rBV2	124393	196171	0.66%	0.135%
30	7.478	712	721	725	rBV2	1799364	3523132	11.84%	2.430%
31	7.531	725	730	733	rBV	2019192	3316485	11.14%	2.288%
32	7.737	753	765	776	rVB	481460	1368799	4.60%	0.944%
33	7.919	789	796	803	rVB2	655166	1228753	4.13%	0.848%
34	8.001	803	810	814	rBV3	117909	239150	0.80%	0.165%
35	8.048	814	818	827	rVB2	174240	292794	0.98%	0.202%
36	8.213	833	846	855	rVV	12339274	29767196	100.00%	20.532%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
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37	8.466	879	889	896	rBV2	243163	571173	1.92%	0.394%
38	8.525	896	899	904	rVB	182492	256334	0.86%	0.177%
39	8.601	904	912	914	rBV2	242003	468323	1.57%	0.323%
40	8.637	914	918	925	rVB	720347	1205821	4.05%	0.832%
41	8.701	925	929	934	rBV	251389	388663	1.31%	0.268%
42	8.754	934	938	944	rVB2	121926	224730	0.75%	0.155%
43	8.937	965	969	977	rVB3	90886	191510	0.64%	0.132%
44	9.042	977	987	990	rBV5	167159	374104	1.26%	0.258%
45	9.089	990	995	1000	rVB	532778	824032	2.77%	0.568%
46	9.172	1000	1009	1014	rBV	1402936	2112664	7.10%	1.457%
47	9.278	1017	1027	1035	rVB4	401188	1234073	4.15%	0.851%
48	9.407	1046	1049	1057	rVB3	68402	169174	0.57%	0.117%
49	9.625	1083	1086	1096	rVB3	103394	228061	0.77%	0.157%
50	9.813	1110	1118	1124	rBV2	498734	880592	2.96%	0.607%
51	10.031	1143	1155	1156	rBV3	278130	717695	2.41%	0.495%
52	10.172	1162	1179	1189	rVB2	1077290	4734677	15.91%	3.266%
53	10.272	1191	1196	1202	rVV2	124311	196879	0.66%	0.136%
54	10.354	1202	1210	1220	rVV2	734982	1880666	6.32%	1.297%
55	10.519	1233	1238	1249	rVV	488706	789036	2.65%	0.544%
56	10.731	1265	1274	1280	rVV2	845669	1850588	6.22%	1.276%
57	10.783	1280	1283	1293	rVB	66955	135071	0.45%	0.093%
58	11.095	1330	1336	1349	rBV	60985	151820	0.51%	0.105%
59	11.995	1481	1489	1493	rVV	70333	173729	0.58%	0.120%
60	12.207	1520	1525	1530	rVV	183591	279909	0.94%	0.193%
61	12.260	1530	1534	1542	rVB	206594	327516	1.10%	0.226%
62	13.389	1720	1726	1733	rVV3	161564	279432	0.94%	0.193%
63	13.971	1818	1825	1831	rBV	1189313	1616899	5.43%	1.115%
64	14.071	1837	1842	1845	rVV	146474	222726	0.75%	0.154%
65	14.113	1845	1849	1862	rVB	145370	237878	0.80%	0.164%
66	14.254	1868	1873	1881	rVB	1301090	1895559	6.37%	1.307%
67	14.560	1917	1925	1931	rBV	1127315	1636669	5.50%	1.129%
68	14.748	1953	1957	1967	rVB	325762	500116	1.68%	0.345%
69	14.877	1973	1979	1988	rVB5	111093	196517	0.66%	0.136%
70	15.548	2085	2093	2100	rBV	1647549	2329476	7.83%	1.607%
71	15.660	2106	2112	2120	rBV	479348	651590	2.19%	0.449%
72	16.236	2204	2210	2214	rBV	111647	156125	0.52%	0.108%
73	16.442	2240	2245	2251	rVB2	77226	140515	0.47%	0.097%
74	17.295	2381	2390	2395	rBV	1200862	1724601	5.79%	1.190%
75	17.395	2401	2407	2417	rVB	1629163	2258864	7.59%	1.558%
76	18.195	2536	2543	2561	rBV	464517	792460	2.66%	0.547%

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77	18.918	2659	2666	2673	rVV8	64085	161049	0.54%	0.111%
78	19.189	2706	2712	2718	rVB4	107565	190832	0.64%	0.132%
79	19.295	2718	2730	2733	rBV2	264206	582052	1.96%	0.401%
80	19.324	2733	2735	2736	rVV	169481	169319	0.57%	0.117%
81	19.348	2736	2739	2747	rVV	404385	609827	2.05%	0.421%
82	19.465	2756	2759	2762	rBV	126440	136672	0.46%	0.094%
83	19.612	2780	2784	2788	rVB	162135	188724	0.63%	0.130%
84	19.677	2790	2795	2800	rBV	1635944	2245589	7.54%	1.549%
85	19.777	2807	2812	2815	rBV3	101575	142141	0.48%	0.098%
86	20.542	2935	2942	2946	rBV3	303368	574018	1.93%	0.396%
87	20.848	2987	2994	2999	rBV7	83837	198481	0.67%	0.137%
88	20.936	3005	3009	3015	rBV5	99073	217812	0.73%	0.150%
89	21.171	3045	3049	3057	rBV	332919	486207	1.63%	0.335%
90	21.459	3093	3098	3105	rBV	1312113	1863693	6.26%	1.285%
91	21.724	3139	3143	3148	rBV3	146110	236903	0.80%	0.163%
92	22.159	3213	3217	3221	rVB3	226942	307258	1.03%	0.212%
93	23.124	3378	3381	3386	rBV2	94073	141259	0.47%	0.097%
94	23.312	3408	3413	3417	rBV	542005	843319	2.83%	0.582%
95	23.383	3417	3425	3434	rVB	7330342	12805981	43.02%	8.833%
96	23.694	3472	3478	3485	rBV	946770	1764275	5.93%	1.217%
97	23.847	3498	3504	3518	rVB2	680569	1441951	4.84%	0.995%
98	24.630	3631	3637	3646	rVB	750608	1757213	5.90%	1.212%
99	24.853	3669	3675	3691	rVB	420106	993337	3.34%	0.685%
100	26.394	3932	3937	3962	rVB3	321659	1092344	3.67%	0.753%

Sum of corrected areas: 144982028

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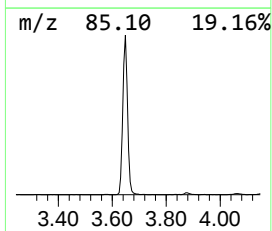
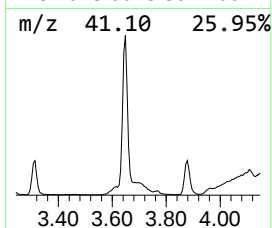
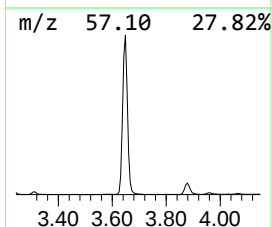
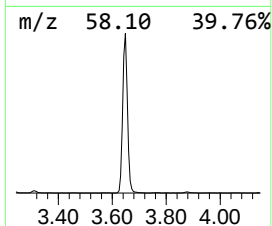
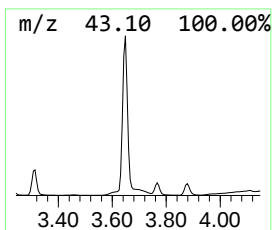
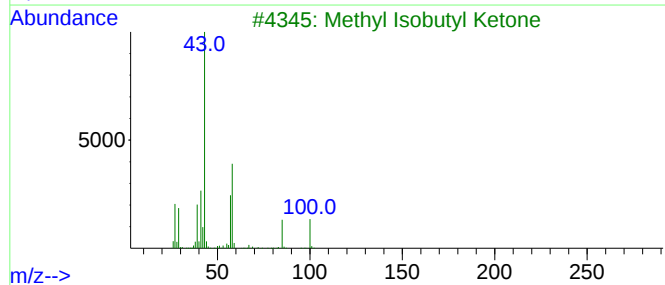
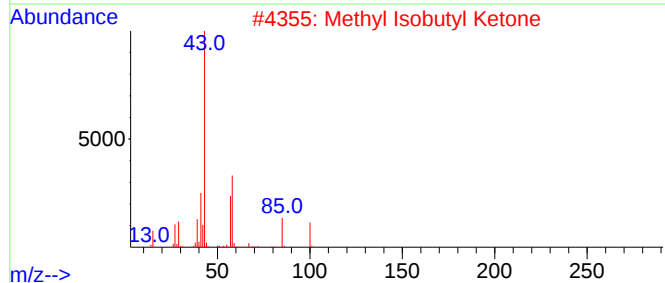
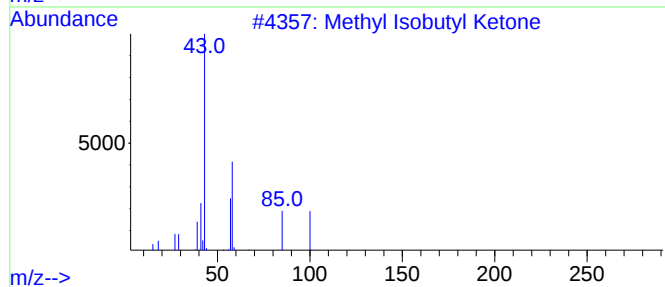
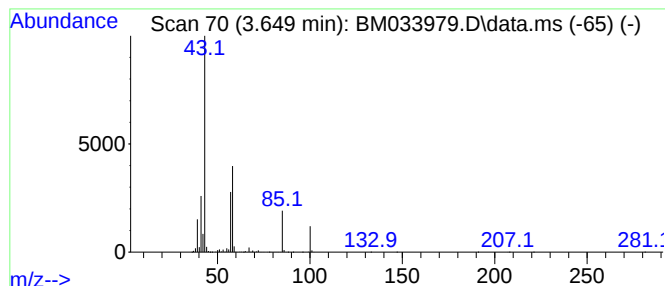
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.649	45.61 ng/ul	2802050	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	91
2			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	87
3			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	83
4			2-Hexanone	100	C6H12O	000591-78-6	72
5			2-Hexanone	100	C6H12O	000591-78-6	72



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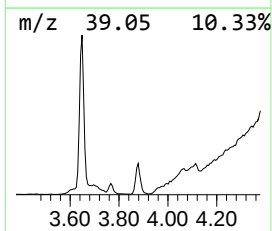
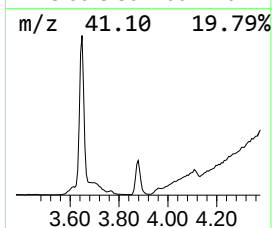
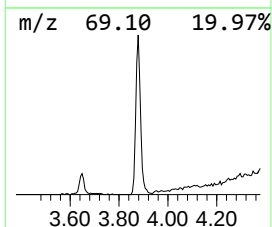
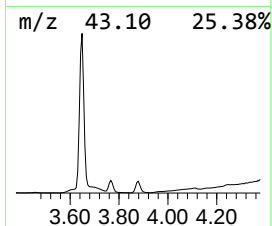
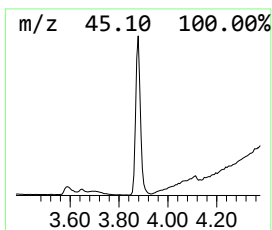
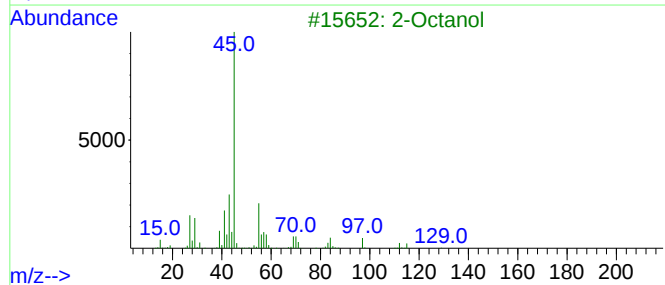
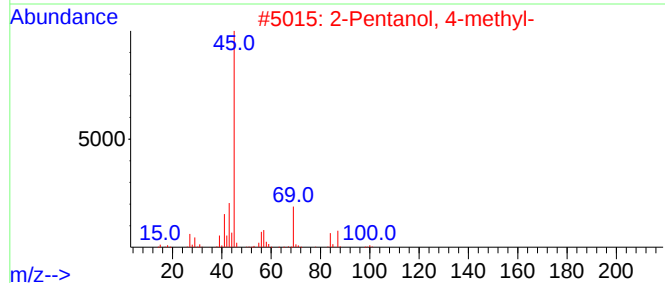
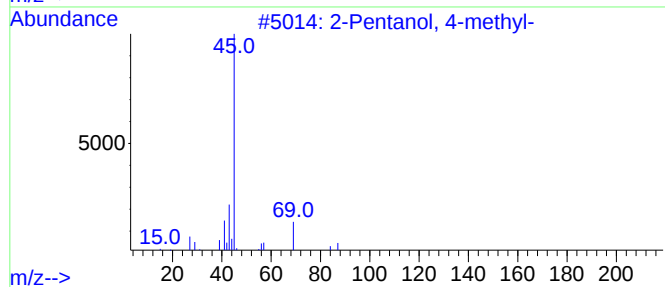
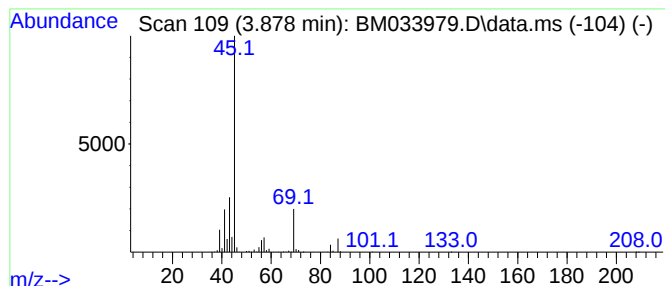
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanol, 4-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.878	10.60 ng/ul	651273	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2		2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
3		2-Octanol	130	C8H18O	000123-96-6	83
4		2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
5		2-Octanol, (R)-	130	C8H18O	005978-70-1	78



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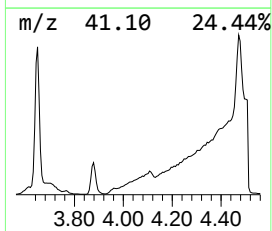
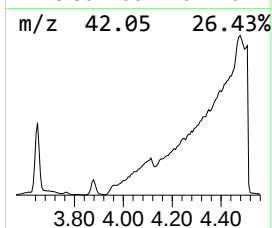
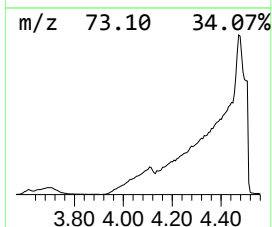
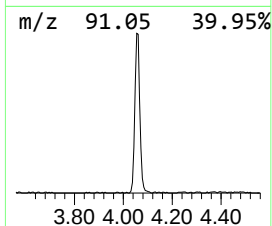
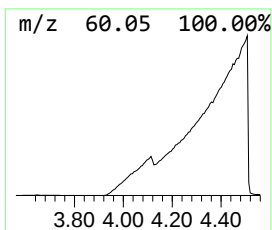
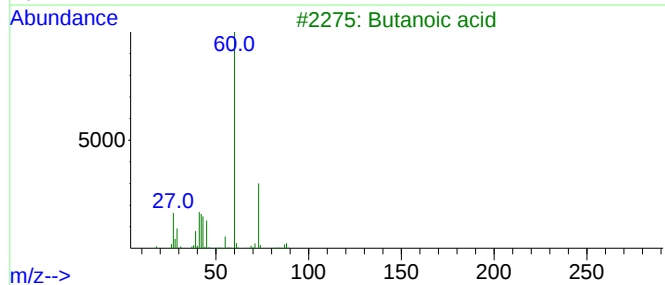
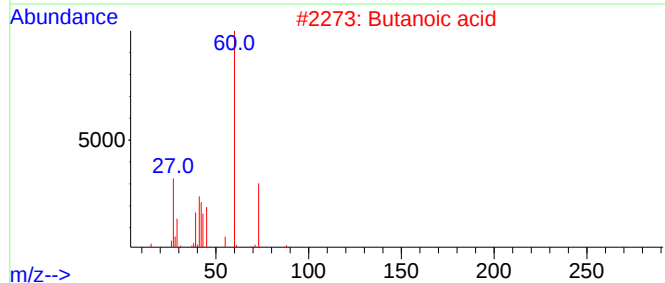
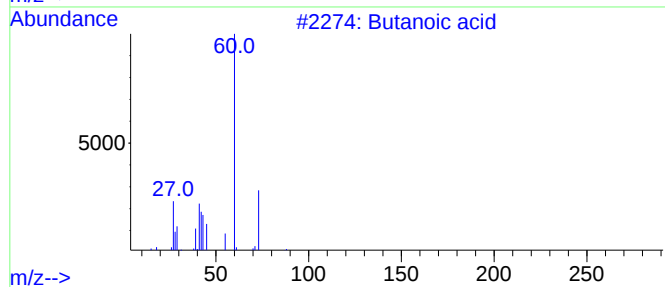
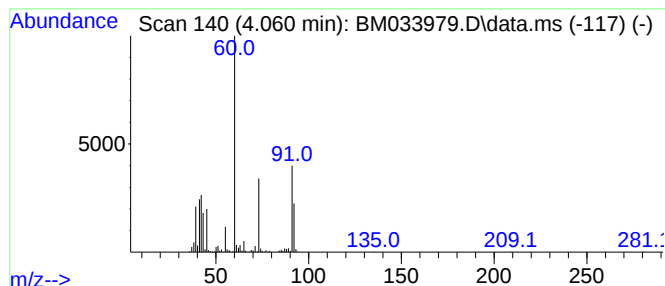
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Butanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.060	22.09 ng/ul	1357340	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid	88	C4H8O2	000107-92-6	68
2			Butanoic acid	88	C4H8O2	000107-92-6	64
3			Butanoic acid	88	C4H8O2	000107-92-6	58
4			Butanoic acid	88	C4H8O2	000107-92-6	53
5			Pentanoic acid	102	C5H10O2	000109-52-4	53



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COVE3

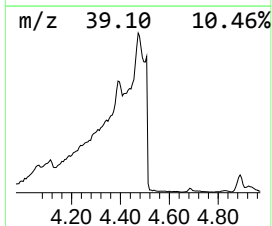
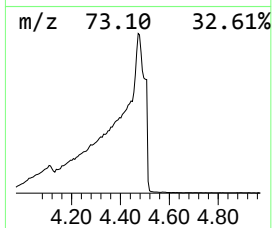
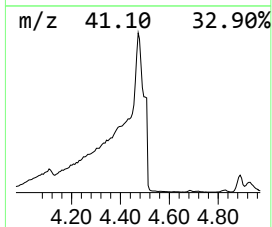
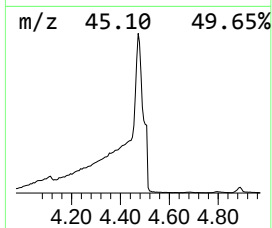
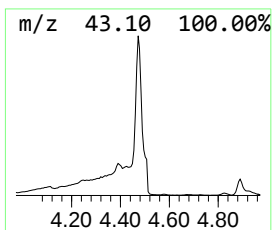
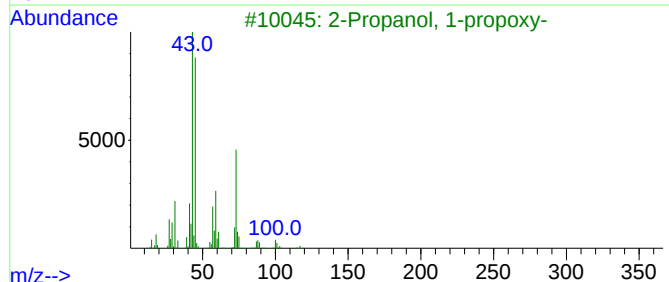
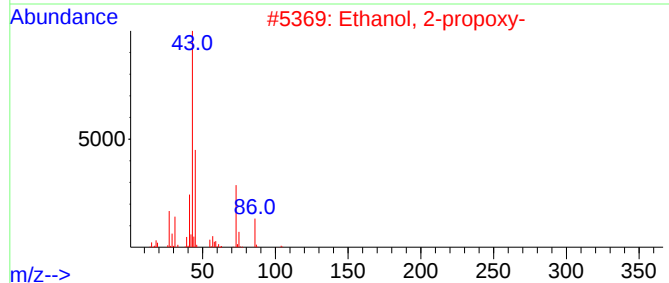
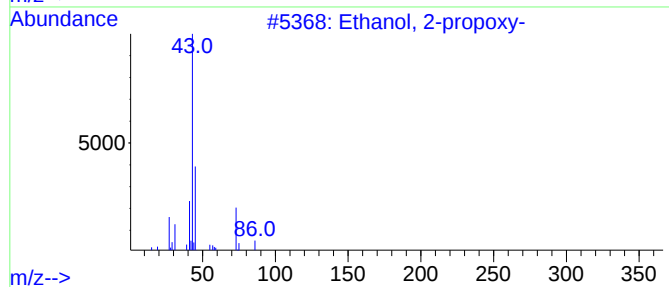
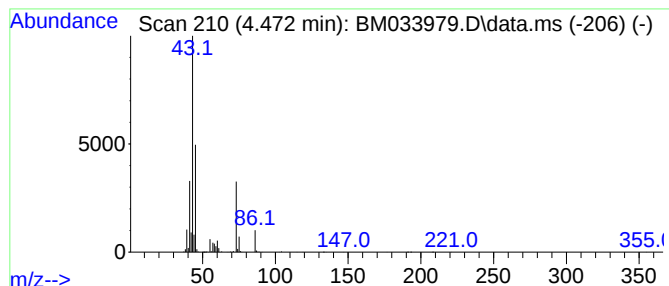
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 2-propoxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.472	25.93 ng/ul	1593050	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-propoxy-	104	C5H12O2	002807-30-9	78
2			Ethanol, 2-propoxy-	104	C5H12O2	002807-30-9	47
3			2-Propanol, 1-propoxy-	118	C6H14O2	001569-01-3	38
4			Ethanol, 2-methoxy-, acetate	118	C5H10O3	000110-49-6	36
5			Ethanol, 2-methoxy-, acetate	118	C5H10O3	000110-49-6	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
COVE3

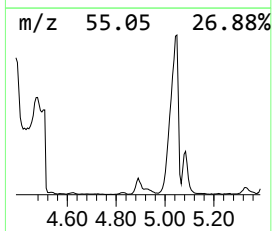
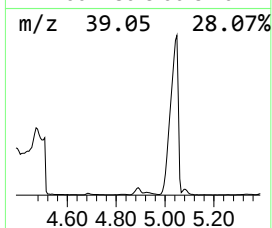
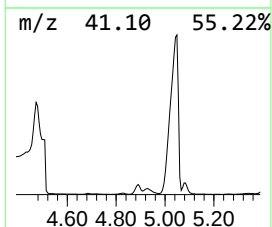
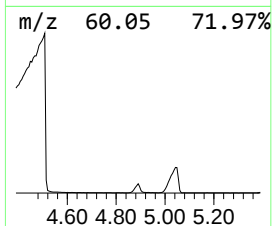
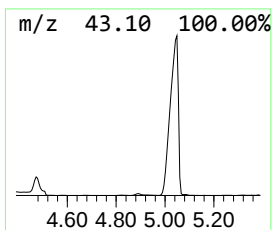
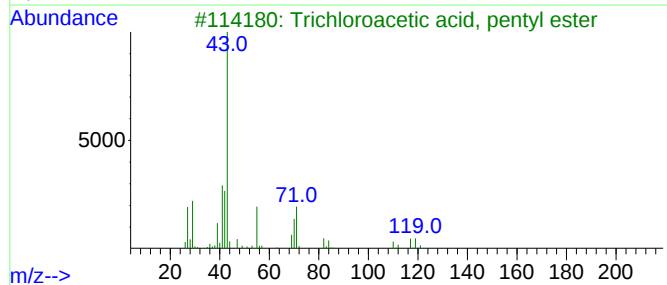
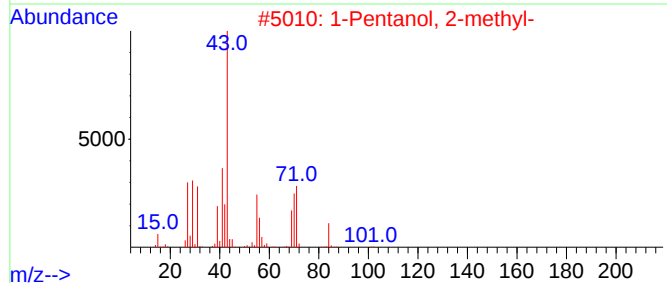
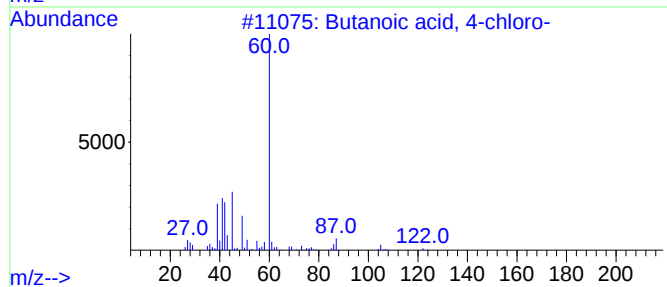
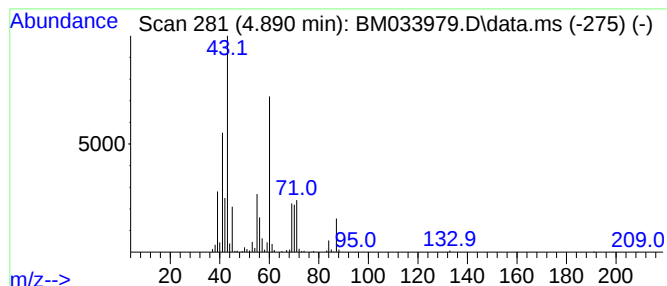
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.890	5.16 ng/ul	316840	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 4-chloro-	122	C4H7ClO2	000627-00-9	43
2			1-Pentanol, 2-methyl-	102	C6H14O	000105-30-6	38
3			Trichloroacetic acid, pentyl ester	232	C7H11Cl3O2	033972-81-5	27
4			1-Pentanol, 2-methyl-	102	C6H14O	000105-30-6	27
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
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Instrument :
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ClientSampleId :
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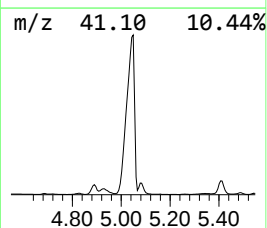
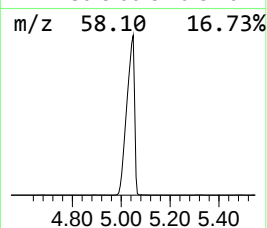
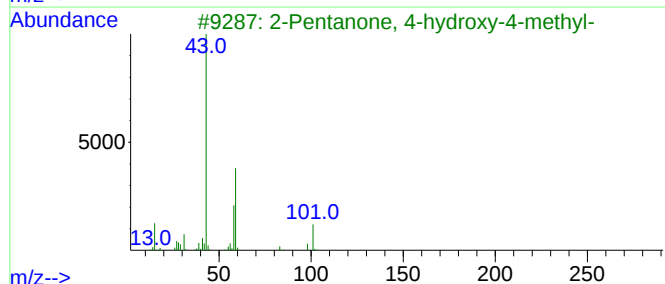
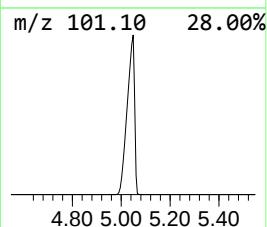
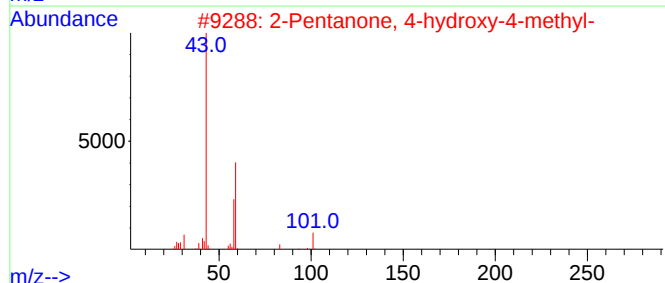
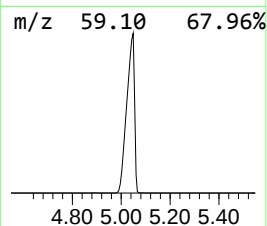
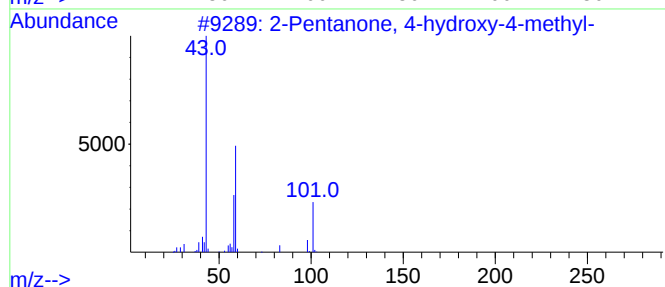
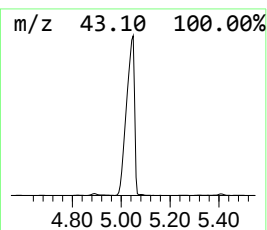
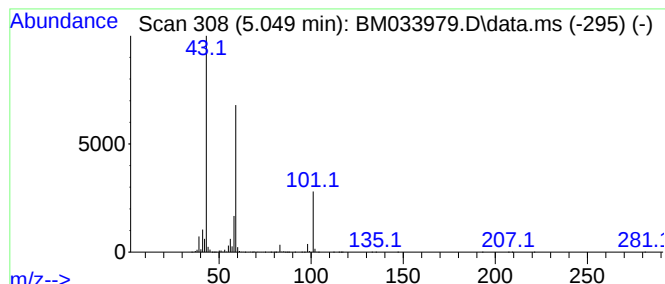
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.049	342.69 ng/ul	21054000	1,4-Dichlorobenzene-d4	7.925

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	42
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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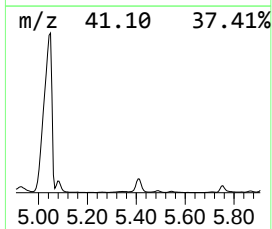
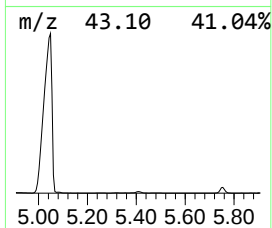
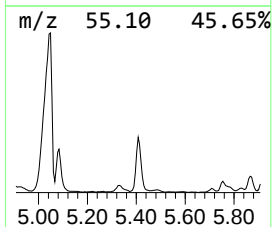
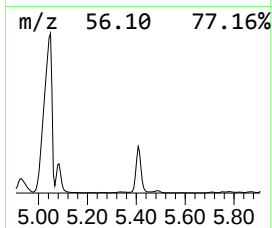
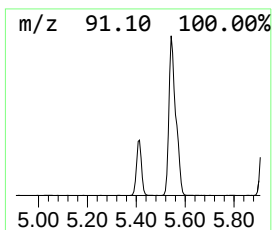
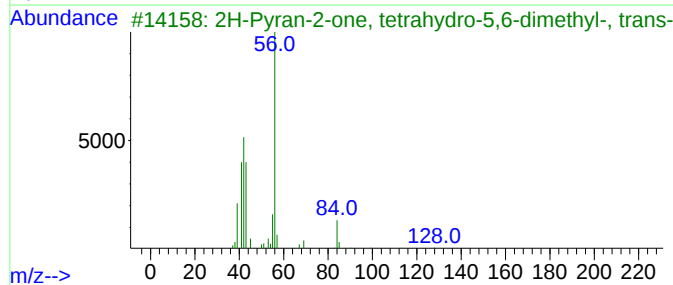
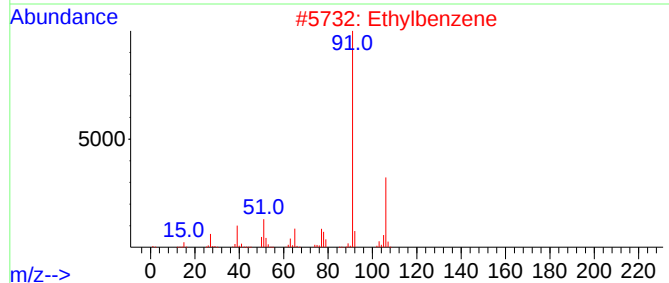
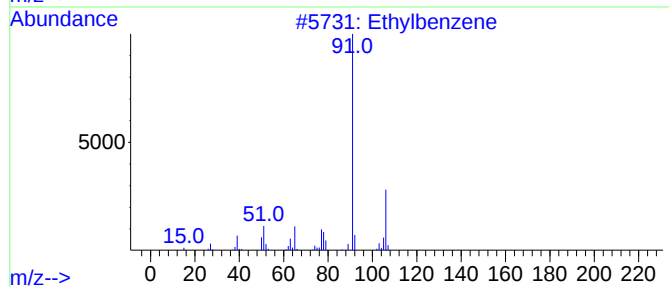
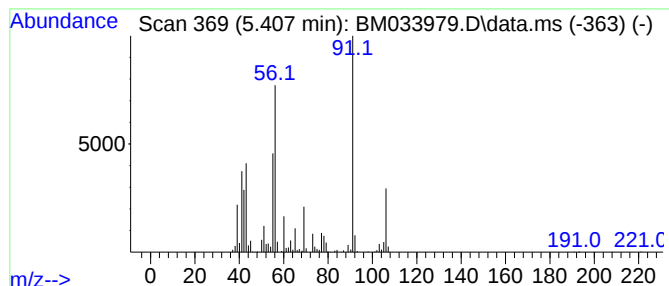
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Ethylbenzene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.407	12.08 ng/ul	742136	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylbenzene	106	C8H10	000100-41-4	64
2			Ethylbenzene	106	C8H10	000100-41-4	60
3			2H-Pyran-2-one, tetrahydro-5,6-d...	128	C7H12O2	024405-16-1	43
4			1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	37
5			Ethylbenzene	106	C8H10	000100-41-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
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Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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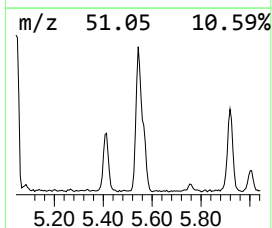
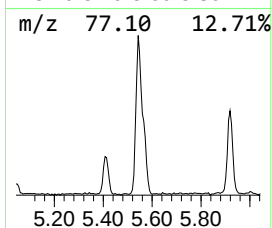
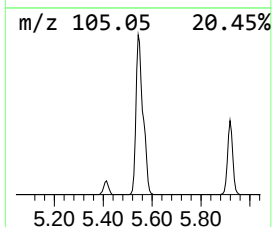
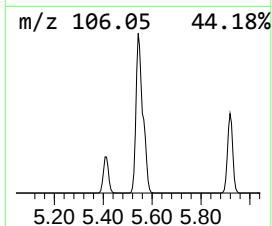
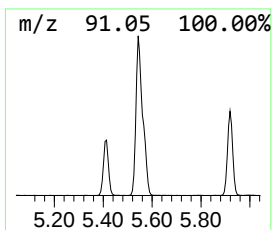
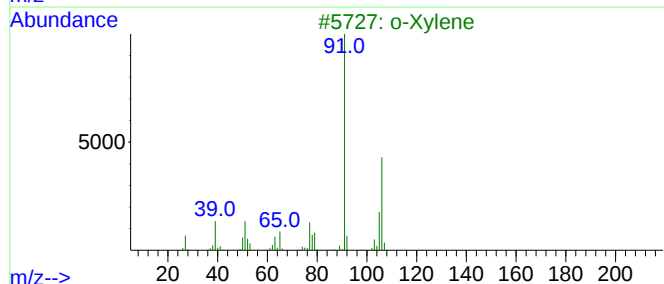
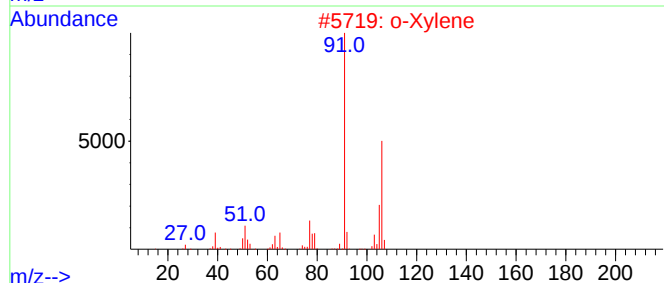
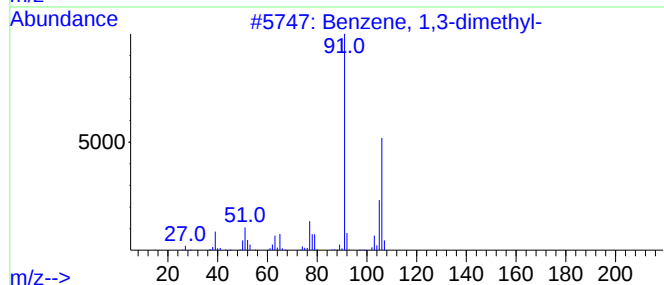
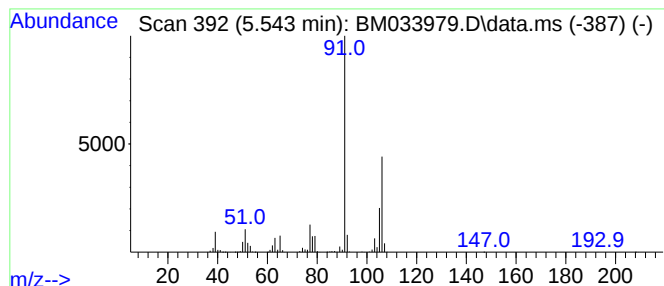
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.543	21.06 ng/ul	1293650	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			o-Xylene	106	C8H10	000095-47-6	97
3			o-Xylene	106	C8H10	000095-47-6	97
4			p-Xylene	106	C8H10	000106-42-3	97
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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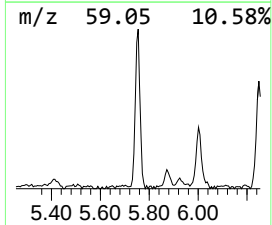
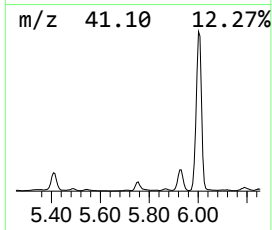
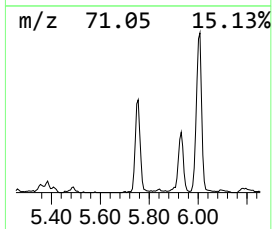
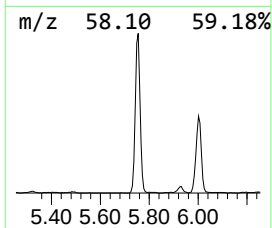
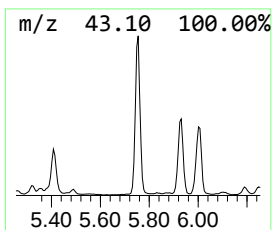
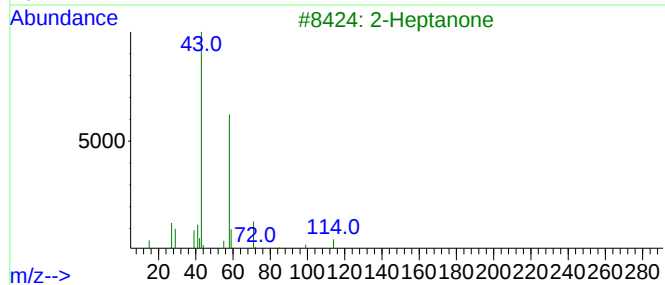
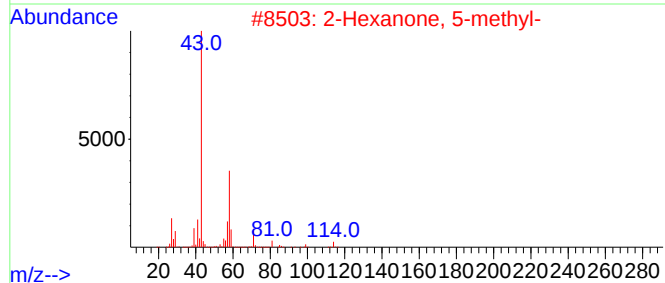
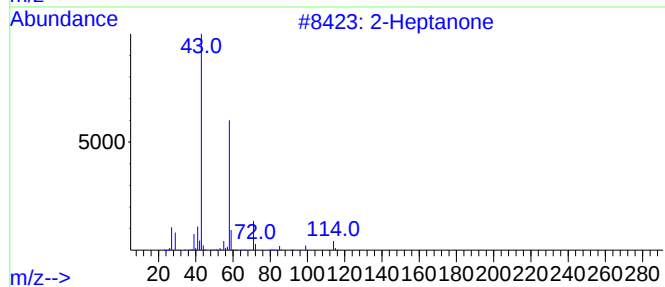
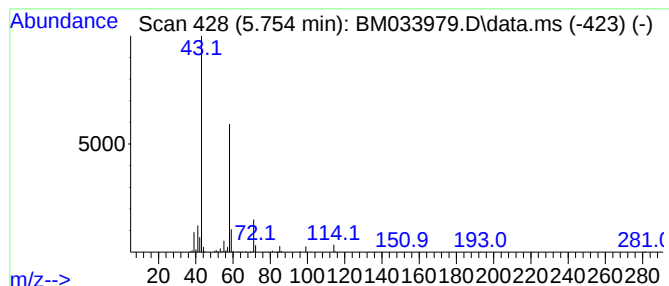
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2-Heptanone Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.754	7.40 ng/ul	454530	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Heptanone	114	C7H14O	000110-43-0	91
2		2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	90
3		2-Heptanone	114	C7H14O	000110-43-0	90
4		2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	86
5		2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
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Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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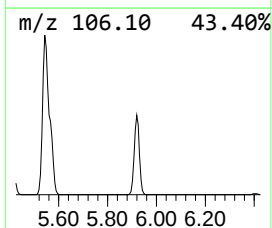
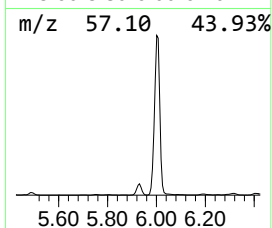
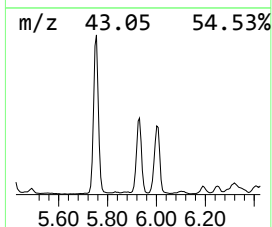
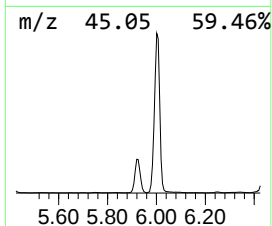
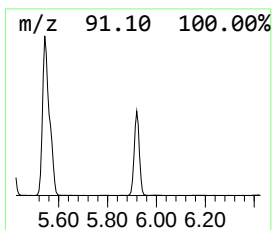
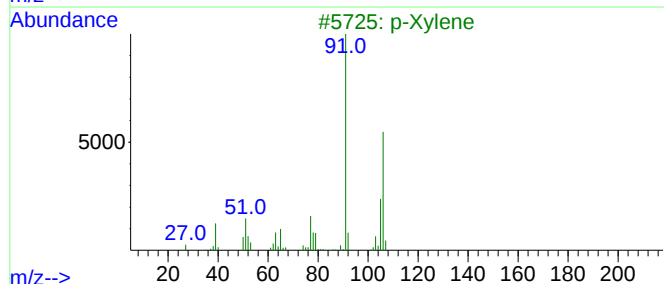
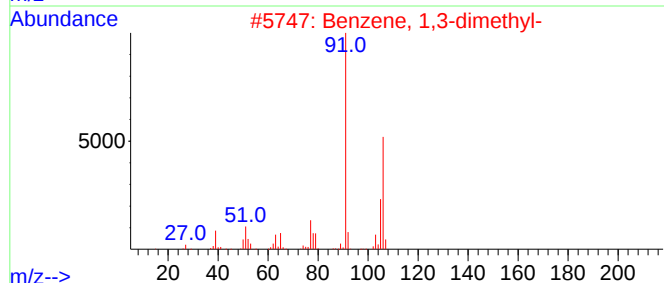
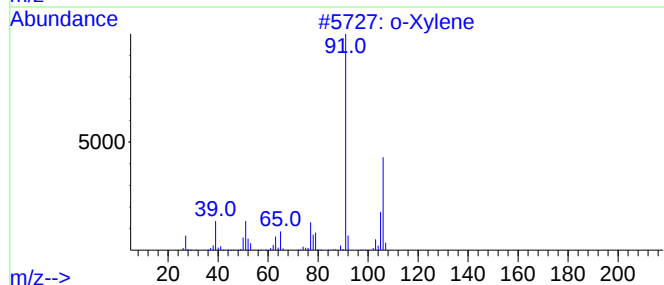
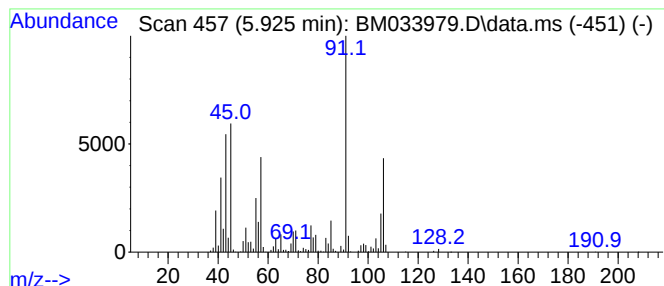
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 o-Xylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.925	17.21 ng/ul	1057160	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	91
2			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	91
3			p-Xylene	106	C8H10	000106-42-3	91
4			p-Xylene	106	C8H10	000106-42-3	91
5			o-Xylene	106	C8H10	000095-47-6	91



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Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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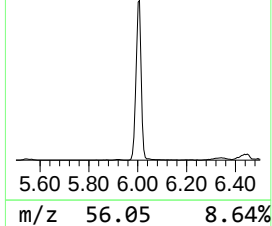
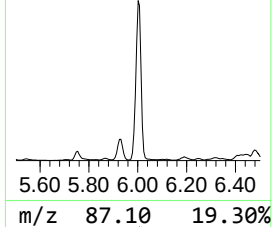
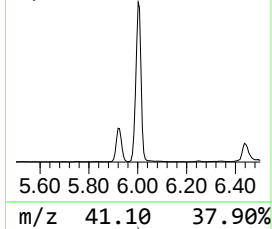
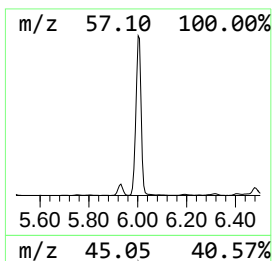
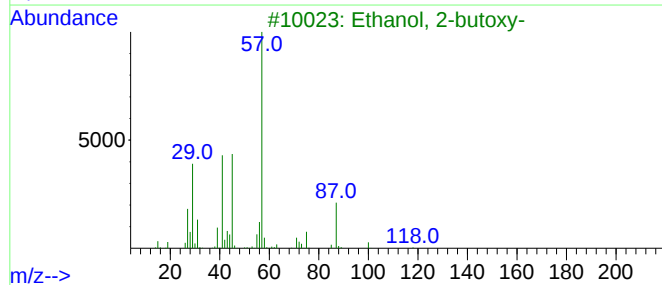
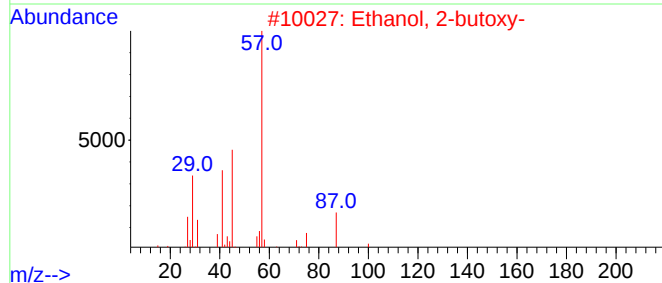
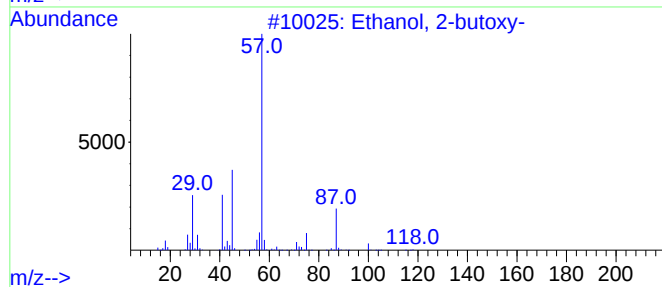
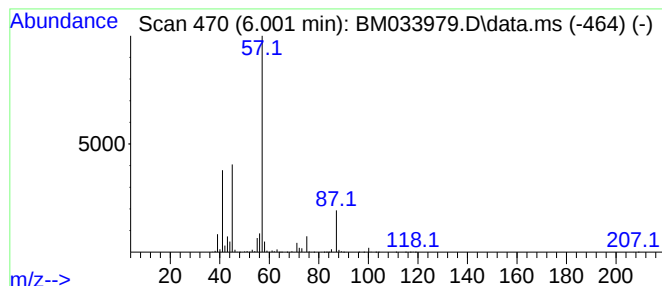
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Ethanol, 2-butoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.001	48.89 ng/ul	3003610	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	91
2			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
3			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
4			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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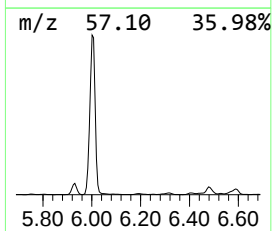
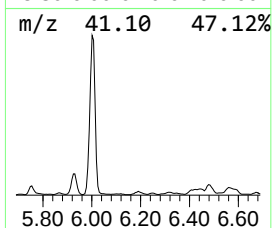
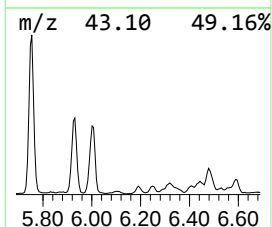
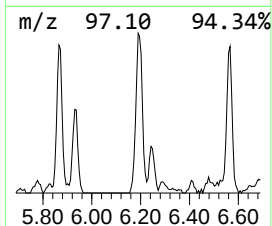
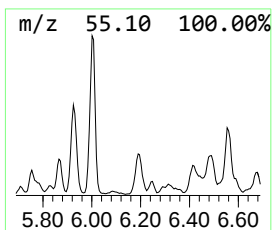
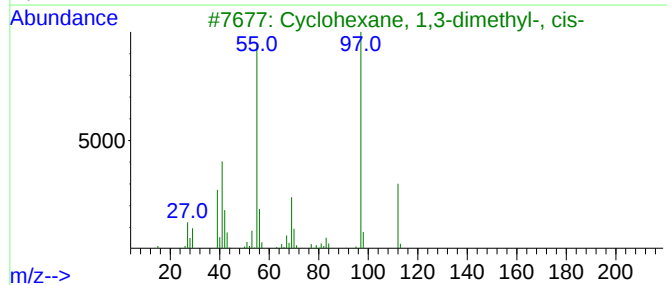
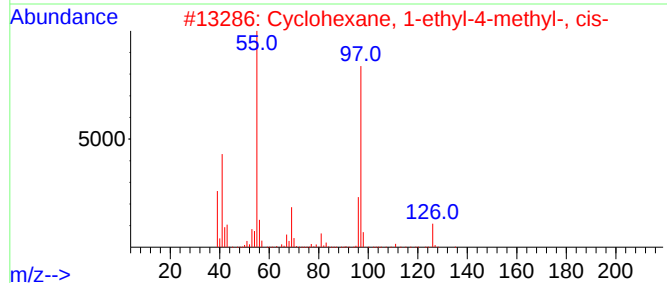
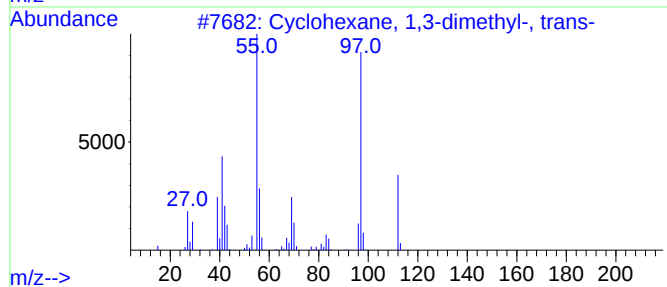
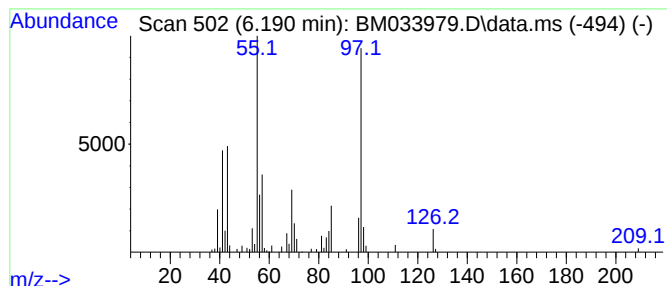
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Cyclic6.190 Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.190	2.36 ng/ul	145094	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	72
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	62
3			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	58
4			Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	58
5			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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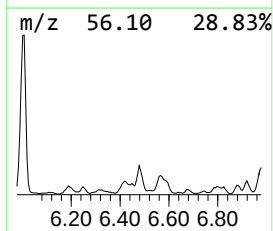
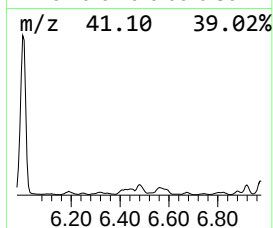
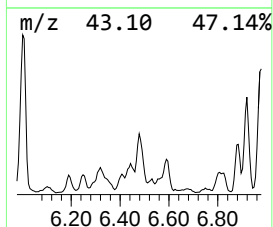
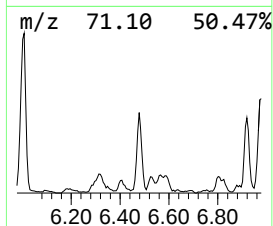
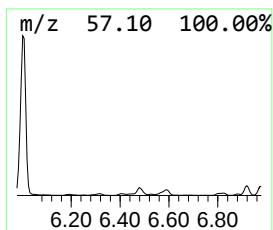
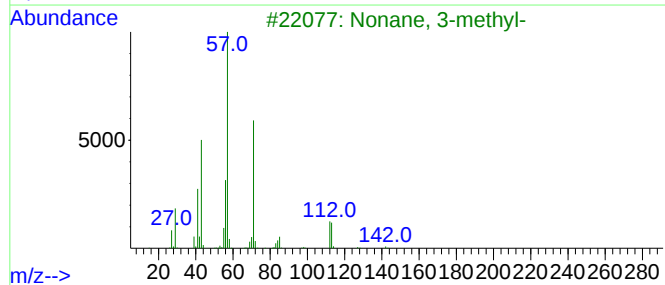
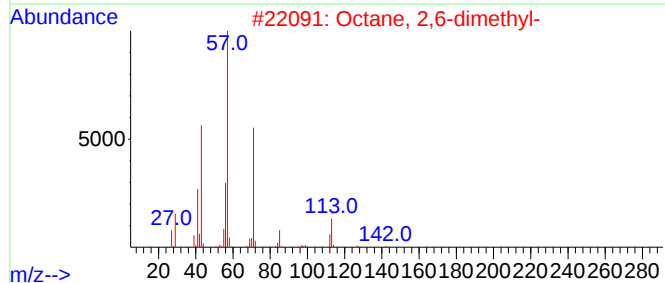
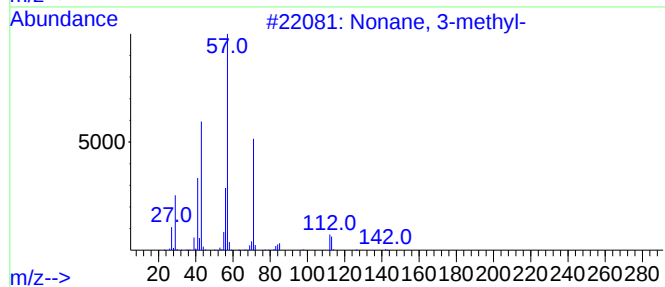
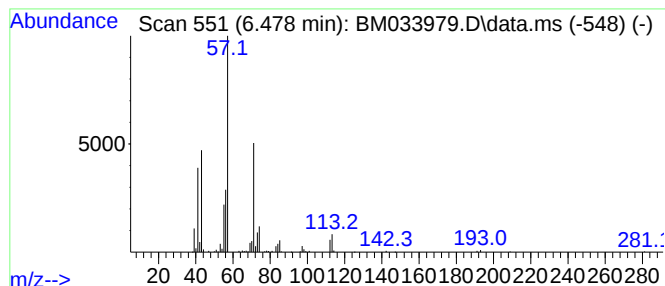
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.478	3.89 ng/ul	238971	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	81
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	64
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
5			Undecane, 3-methyl-	170	C12H26	001002-43-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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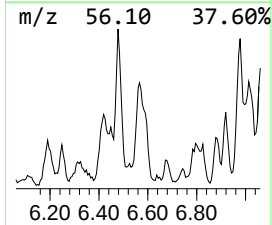
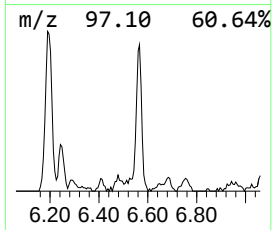
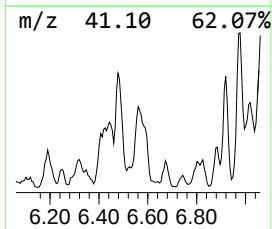
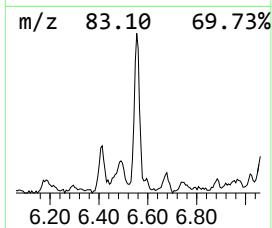
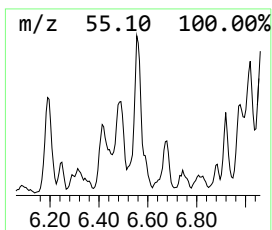
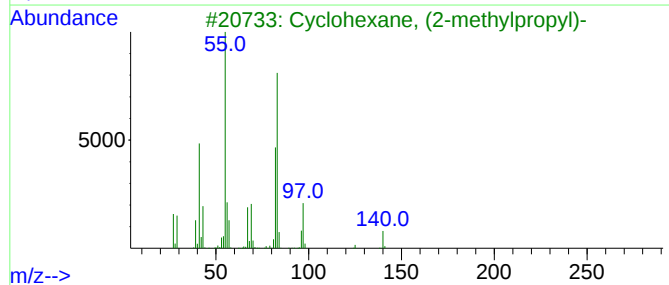
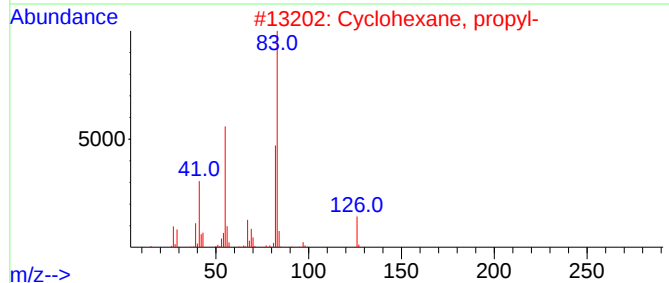
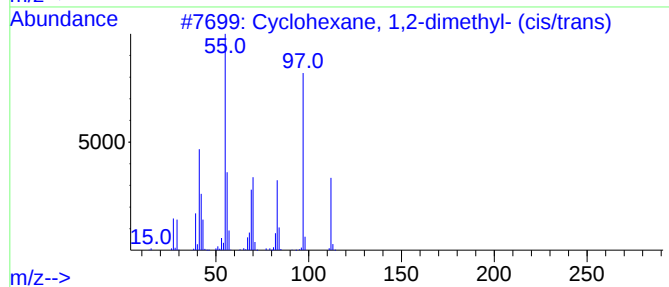
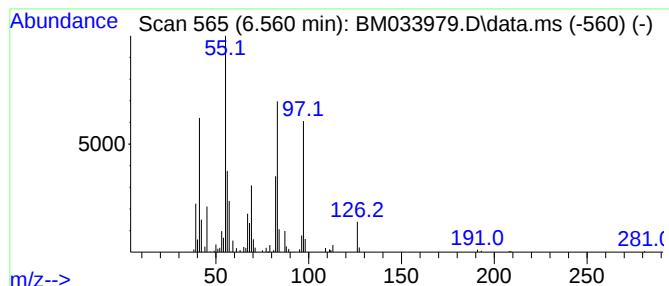
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Cyclic6.560 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.560	2.99 ng/ul	183947	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	47
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	43
3			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	43
4			3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	42
5			2,5-Dimethyl-1-pyrroline	97	C6H11N	000822-64-0	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
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Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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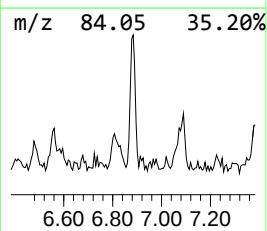
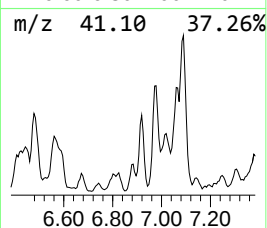
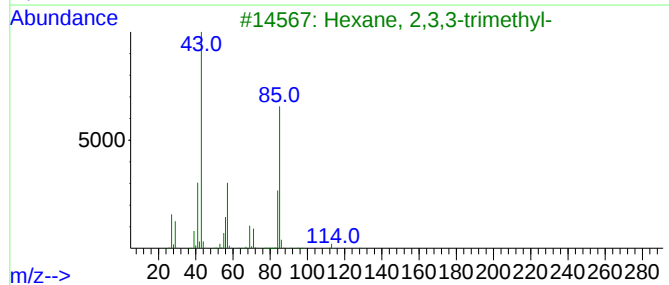
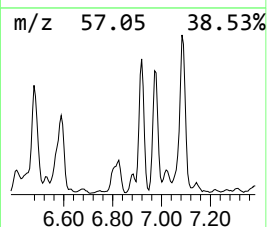
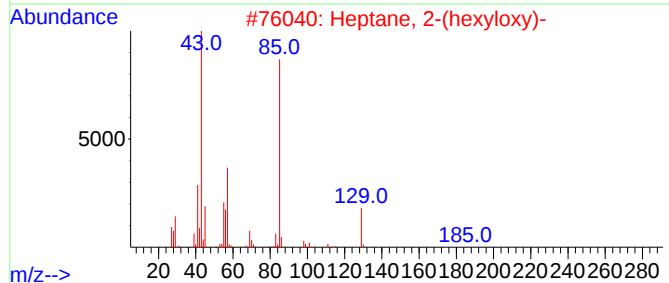
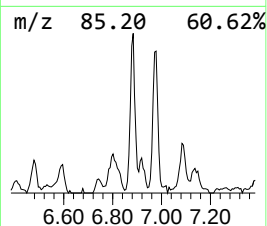
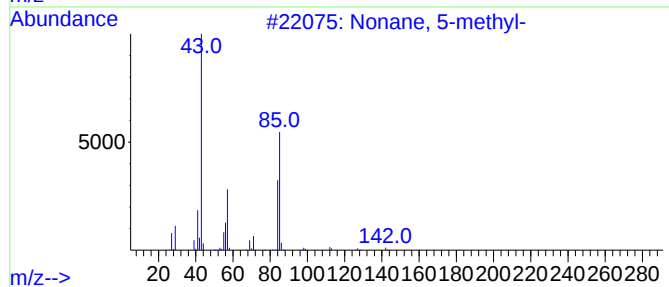
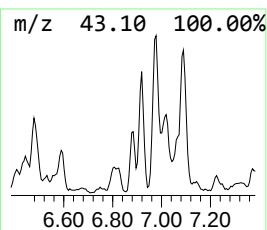
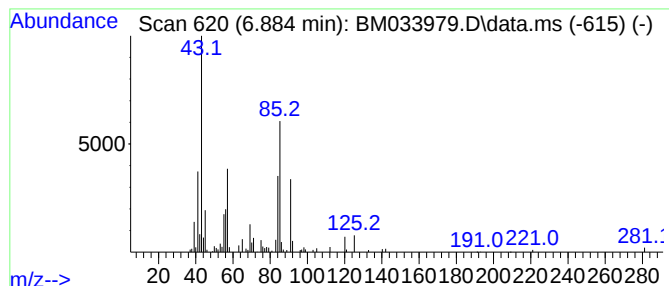
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.884	2.15 ng/ul	132077	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 5-methyl-	142	C10H22	015869-85-9	60
2		Heptane, 2-(hexyloxy)-	200	C13H28O	074663-79-9	53
3		Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	53
4		Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	53
5		Heptane, 3-methyl-	114	C8H18	000589-81-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
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Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
COVE3

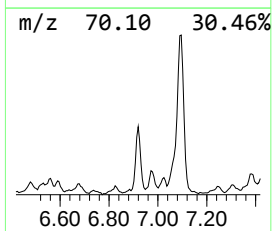
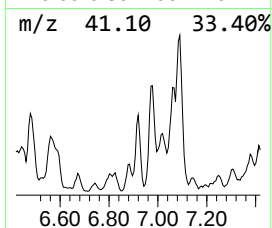
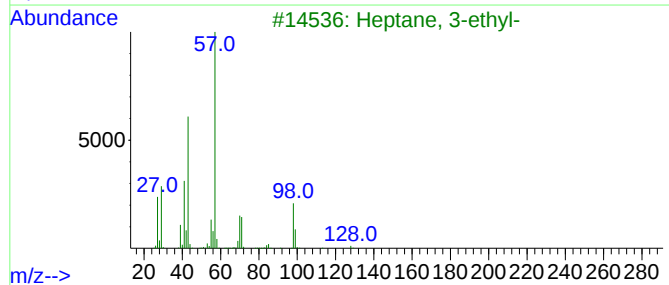
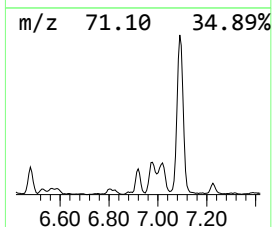
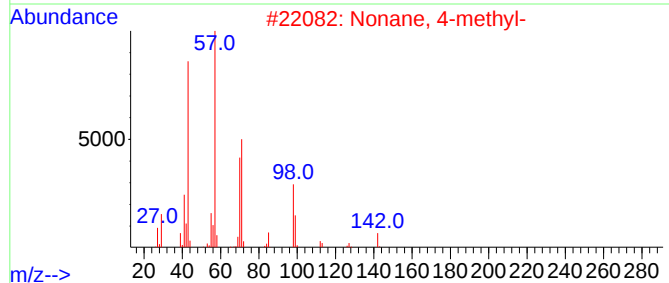
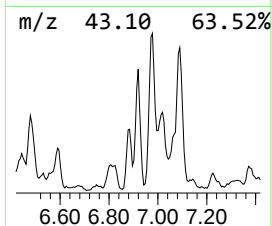
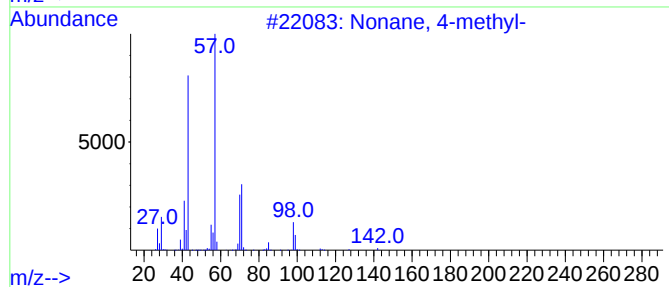
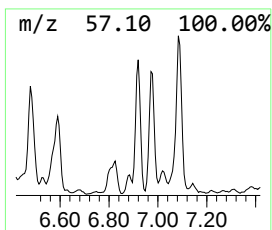
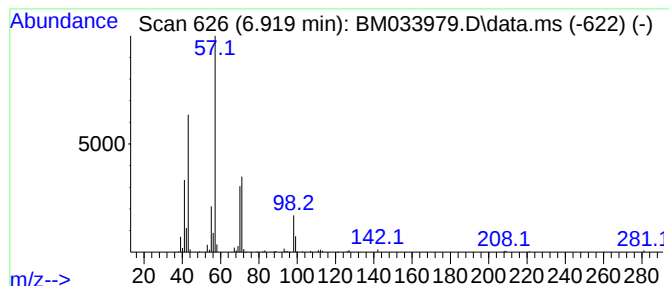
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.919	4.18 ng/ul	257114	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-	142	C10H22	017301-94-9	86	
2	Nonane, 4-methyl-	142	C10H22	017301-94-9	64	
3	Heptane, 3-ethyl-	128	C9H20	015869-80-4	50	
4	Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	50	
5	Methylamine, N-(1-ethylpentylide...	127	C8H17N	018641-73-1	43	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

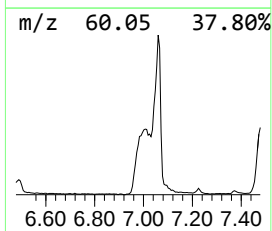
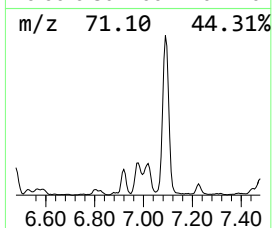
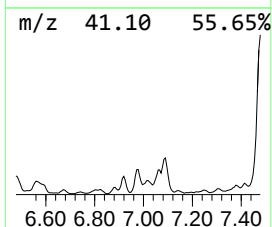
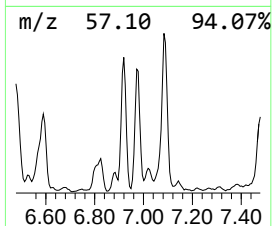
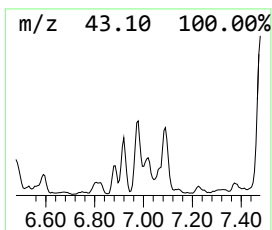
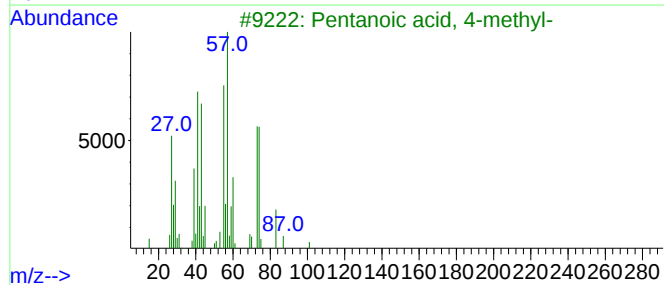
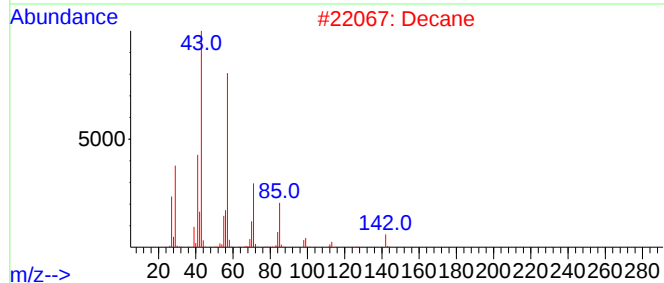
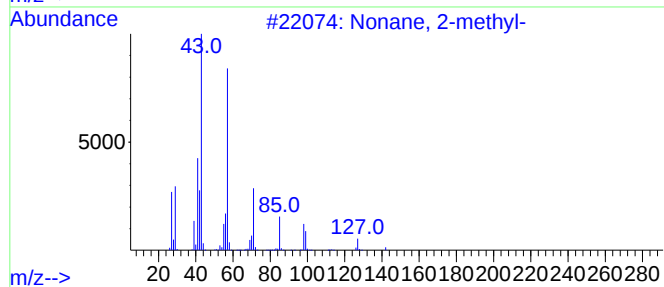
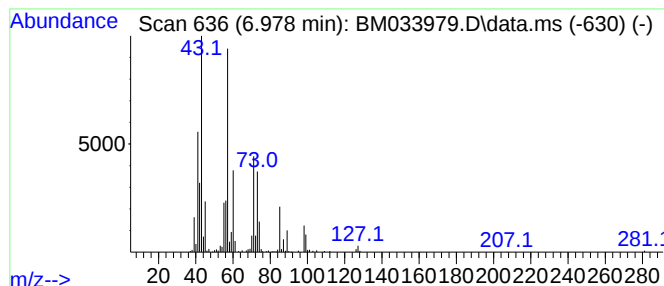
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.978	6.92 ng/ul	425255	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2-methyl-	142	C10H22	000871-83-0	43
2		Decane	142	C10H22	000124-18-5	38
3		Pentanoic acid, 4-methyl-	116	C6H12O2	000646-07-1	38
4		Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	38
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

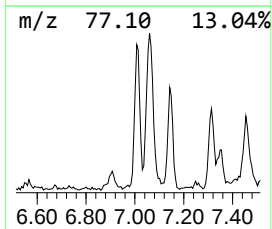
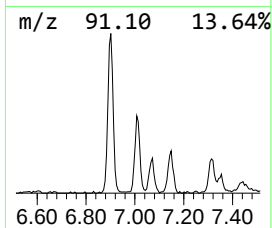
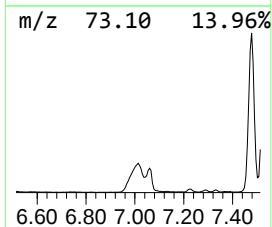
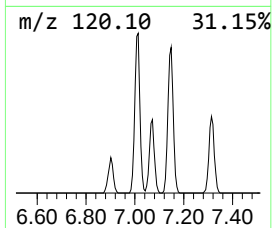
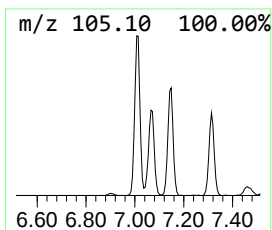
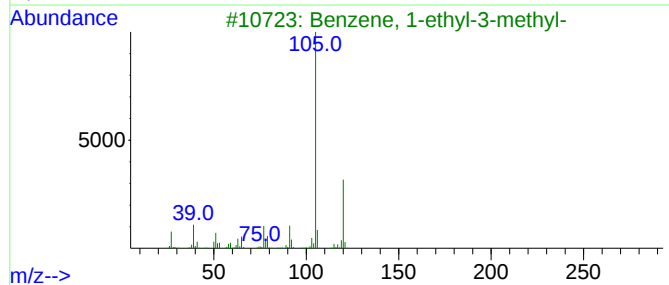
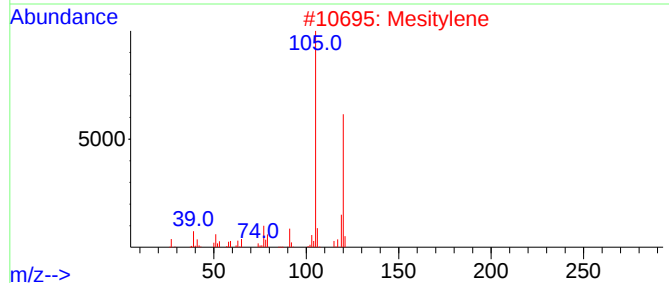
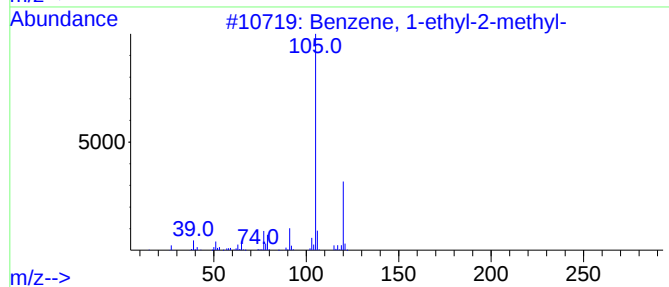
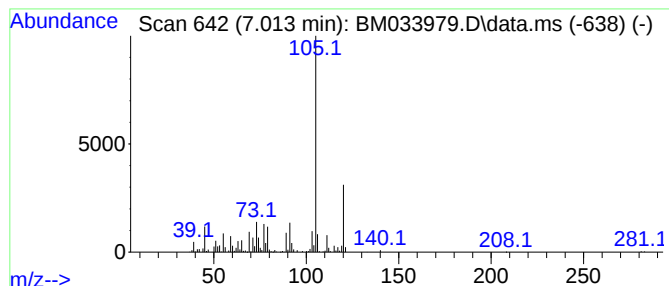
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 1-ethyl-2-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.013	7.66 ng/ul	470732	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	81
2			Mesitylene	120	C9H12	000108-67-8	81
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	81
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	81
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
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Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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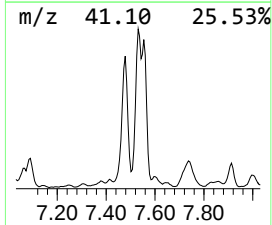
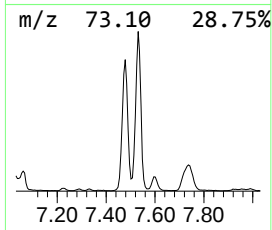
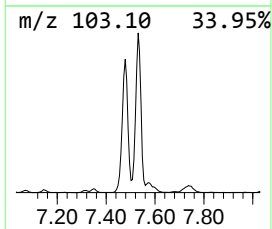
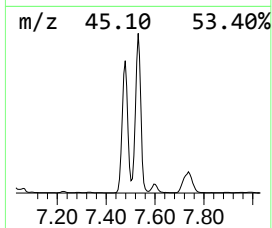
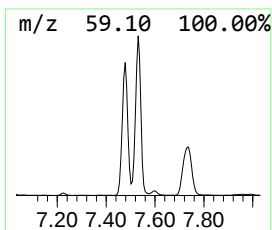
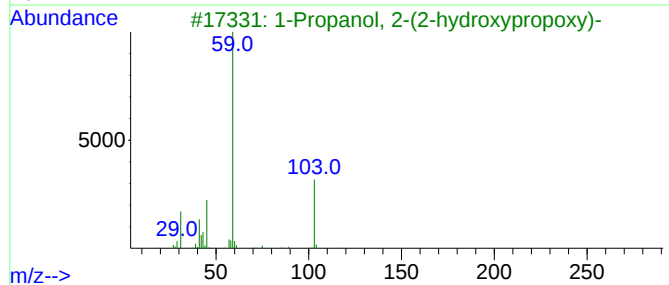
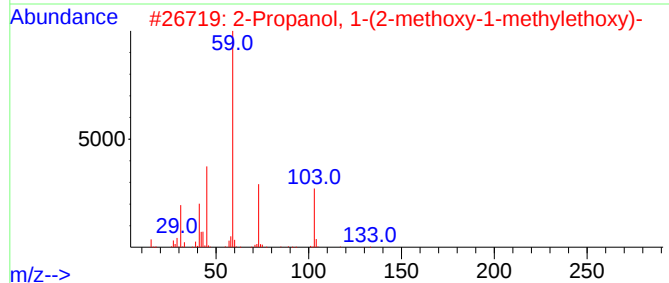
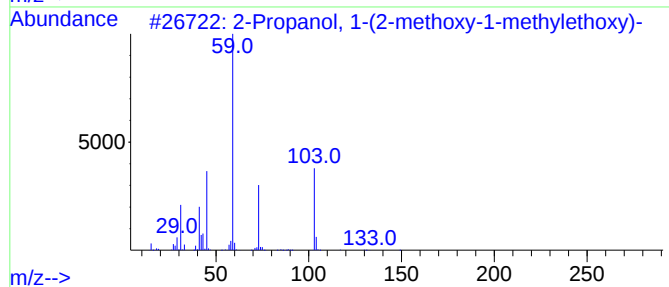
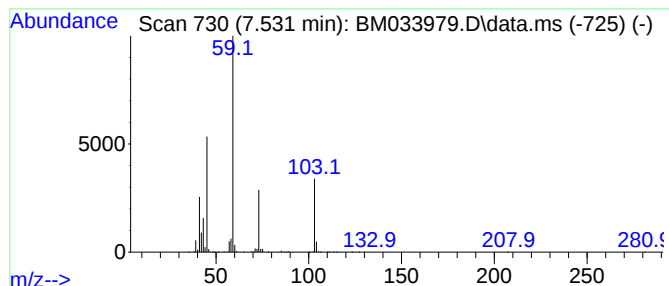
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.531	53.98 ng/ul	3316490	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	86		
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59		
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

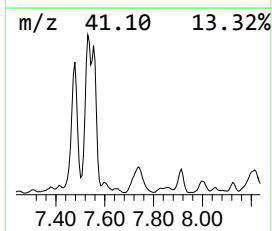
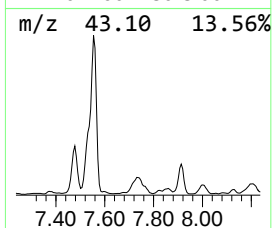
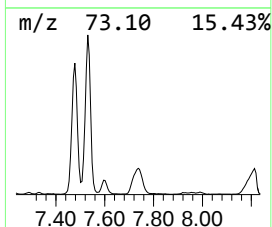
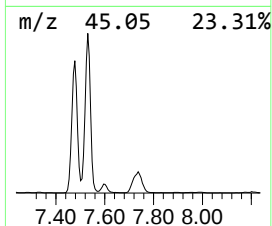
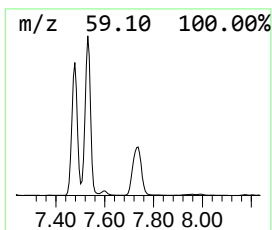
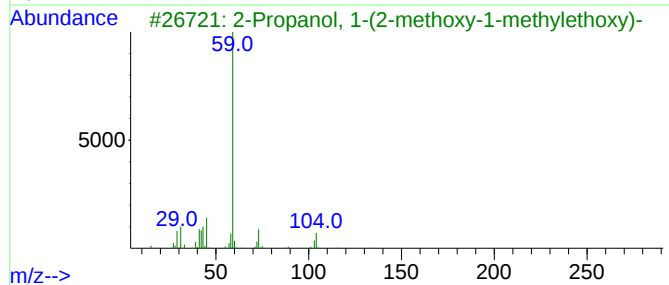
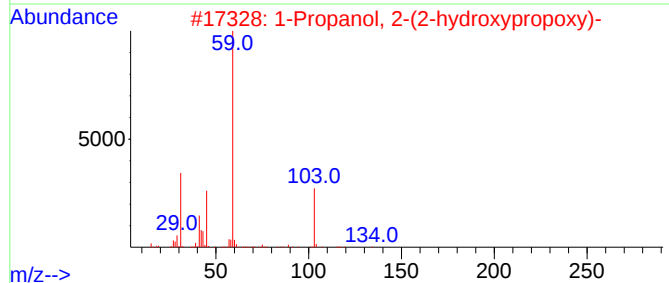
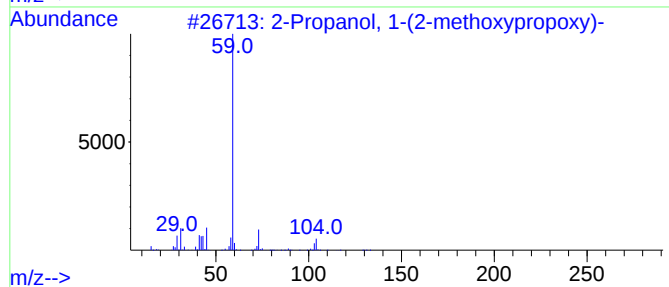
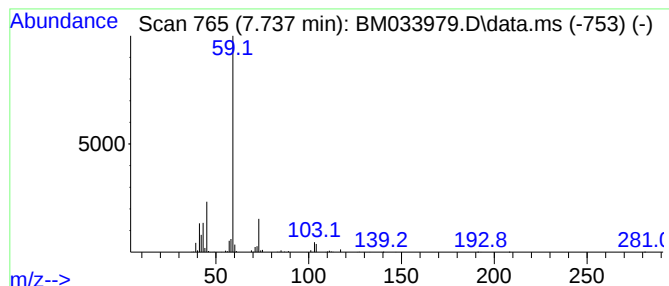
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 2-Propanol, 1-(2-methoxypro... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.737	22.28 ng/ul	1368800	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
2		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
3		2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
4		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
5		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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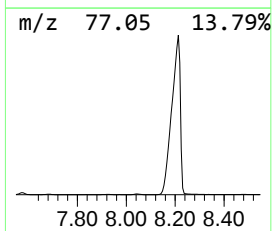
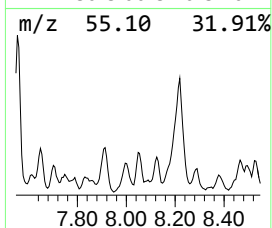
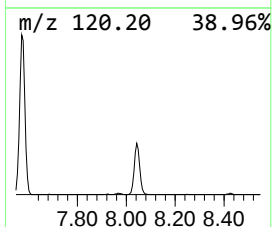
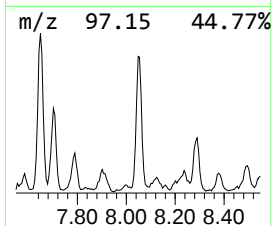
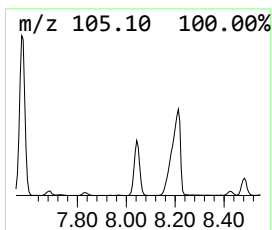
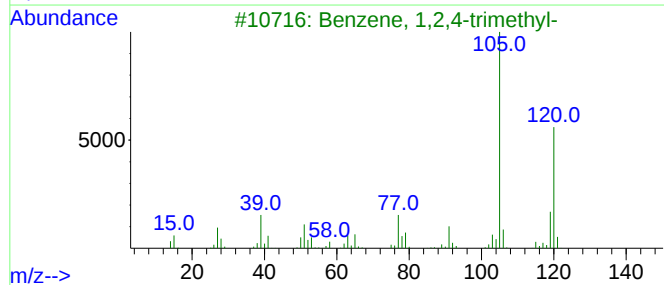
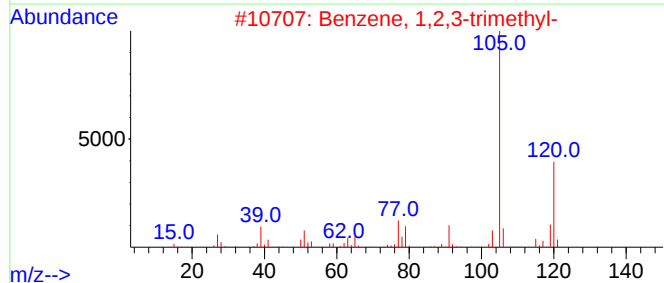
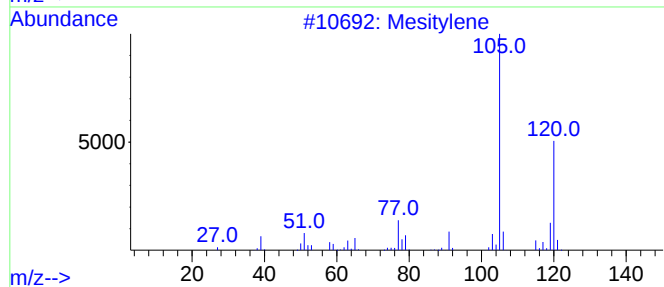
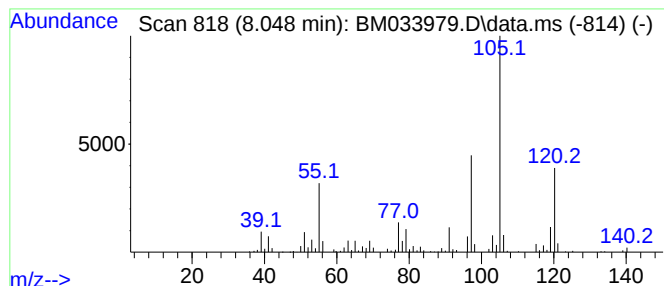
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Mesitylene Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.048	4.77 ng/ul	292794	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	93
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	92
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

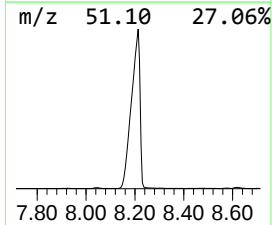
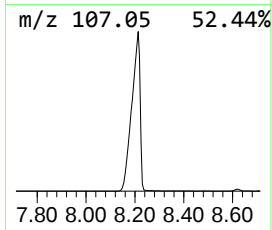
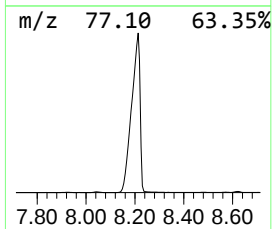
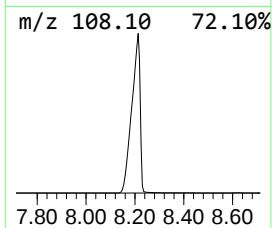
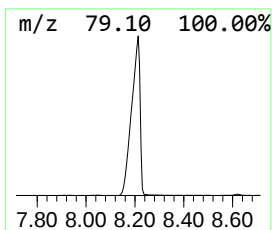
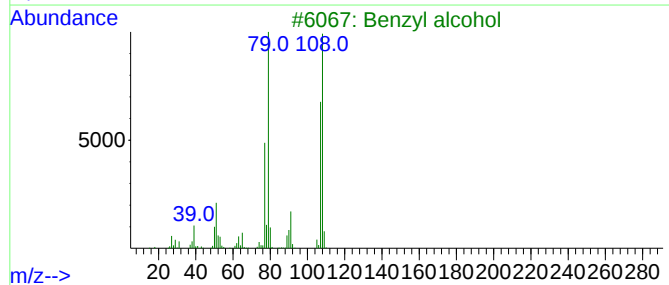
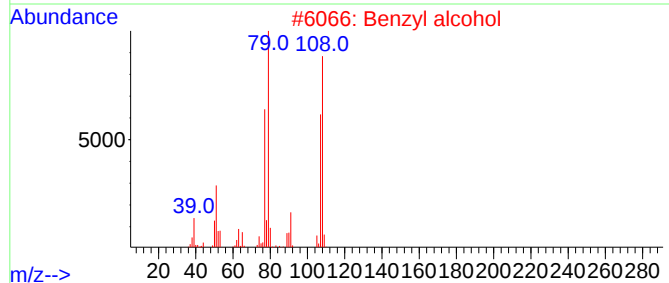
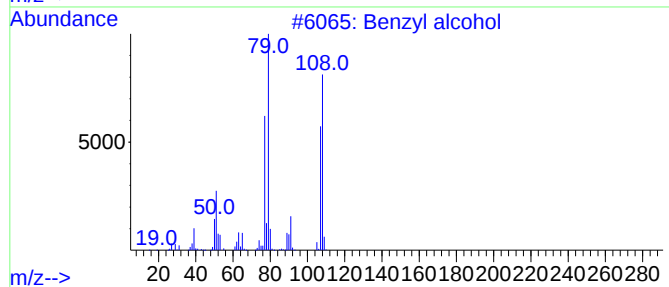
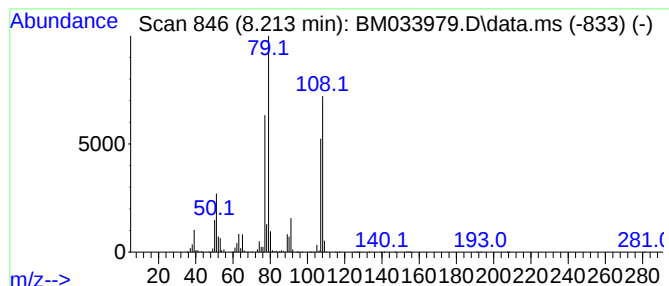
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Benzyl alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.213	484.51 ng/ul	29767200	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	98
2			Benzyl alcohol	108	C7H8O	000100-51-6	97
3			Benzyl alcohol	108	C7H8O	000100-51-6	97
4			Benzyl alcohol	108	C7H8O	000100-51-6	94
5			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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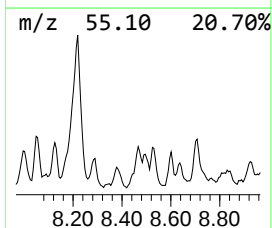
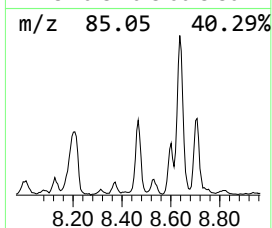
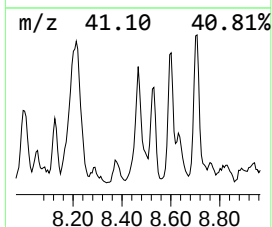
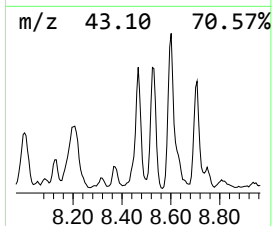
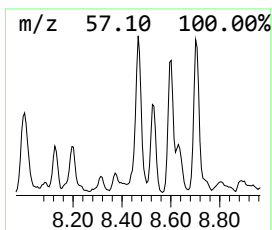
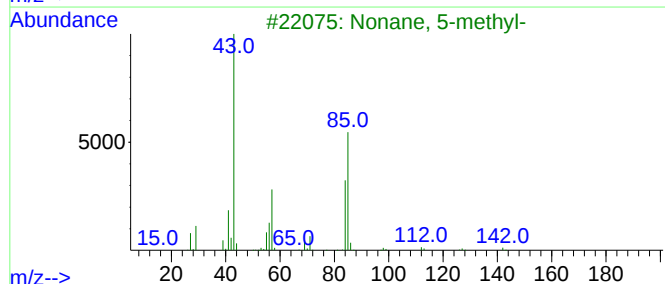
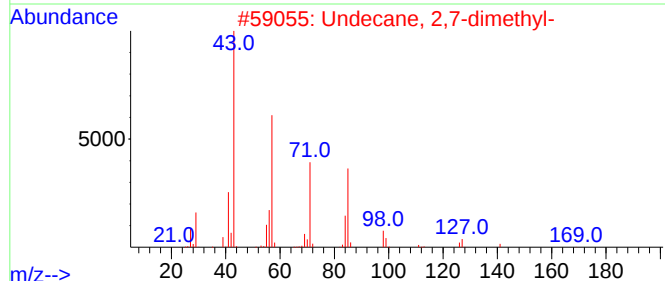
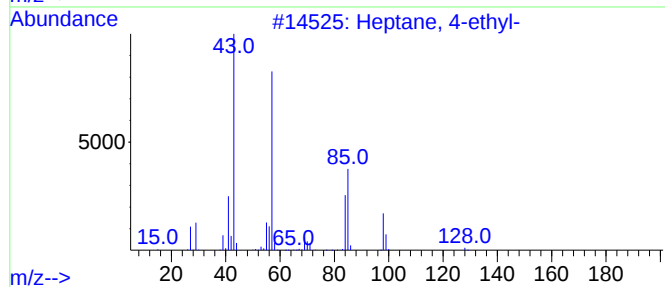
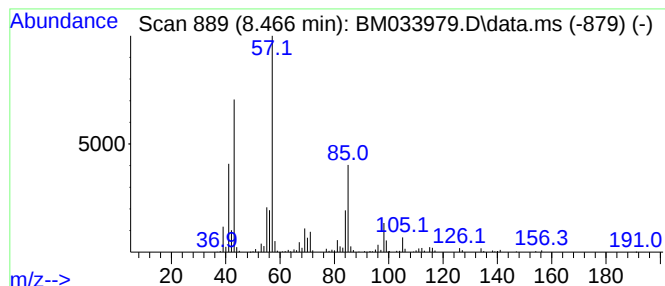
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.466	9.30 ng/ul	571173	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-ethyl-	128	C9H20	002216-32-2	80
2		Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	64
3		Nonane, 5-methyl-	142	C10H22	015869-85-9	59
4		Butanoic acid, 3-oxo-, 1,1-dimet...	158	C8H14O3	001694-31-1	53
5		Dodecane, 5-methyl-	184	C13H28	017453-93-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

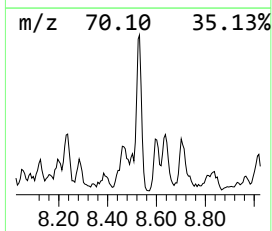
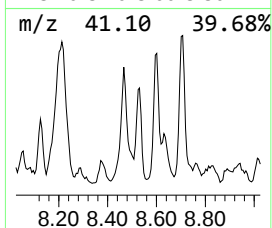
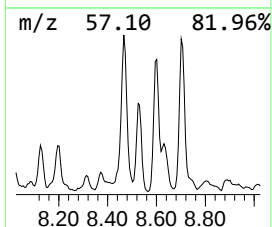
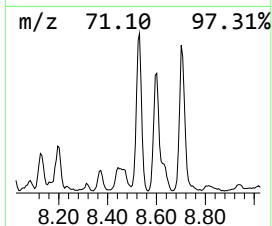
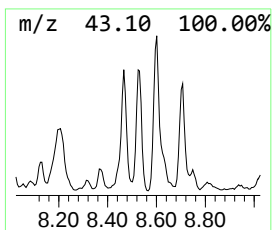
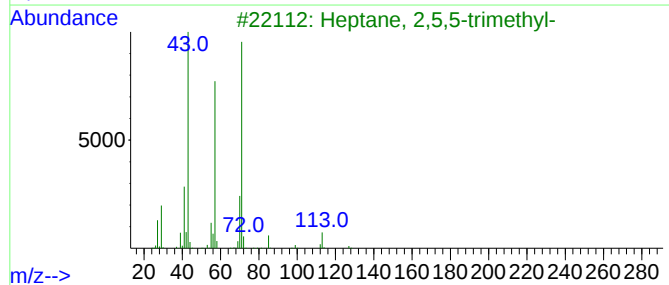
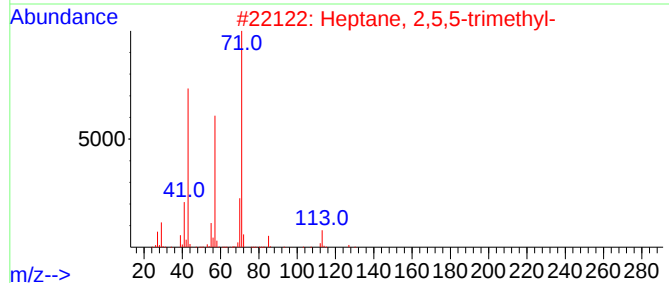
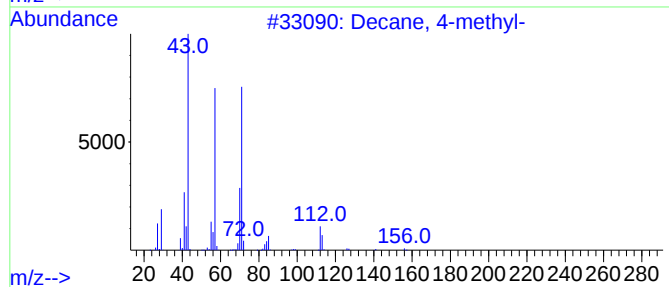
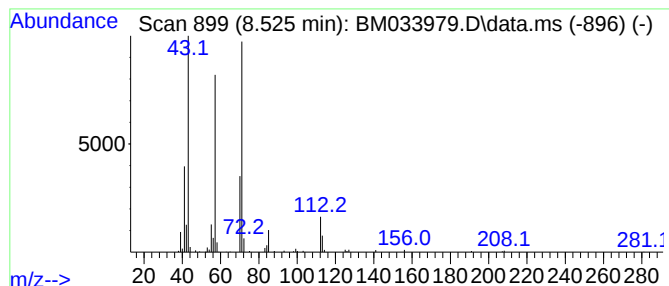
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ClientSampleId :
COVE3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.525	4.17 ng/ul	256334	1,4-Dichlorobenzene-d4			7.925
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-		156	C11H24	002847-72-5	91
2	Heptane, 2,5,5-trimethyl-		142	C10H22	001189-99-7	83
3	Heptane, 2,5,5-trimethyl-		142	C10H22	001189-99-7	78
4	Octane, 3-ethyl-		142	C10H22	005881-17-4	72
5	Tridecane, 4-methyl-		198	C14H30	026730-12-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

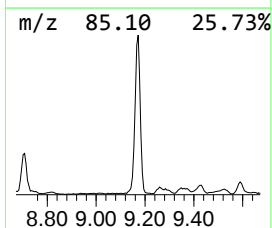
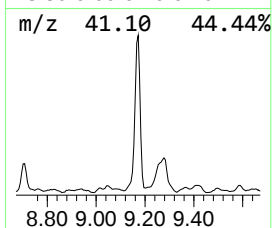
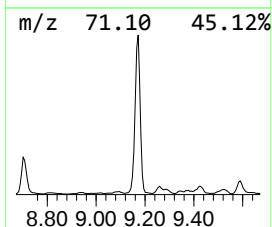
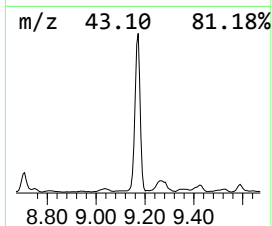
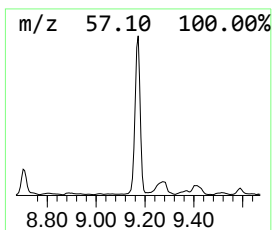
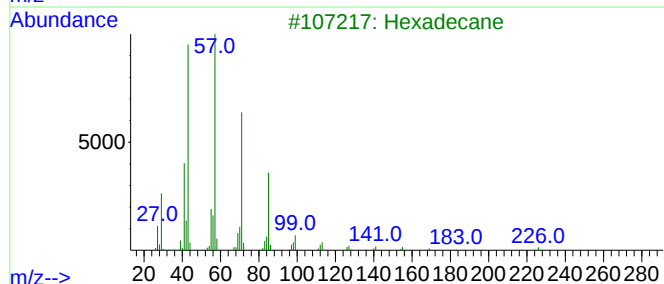
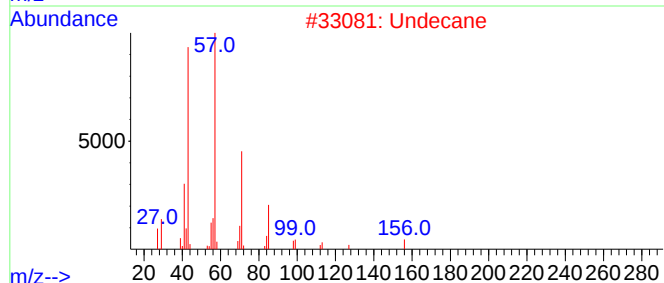
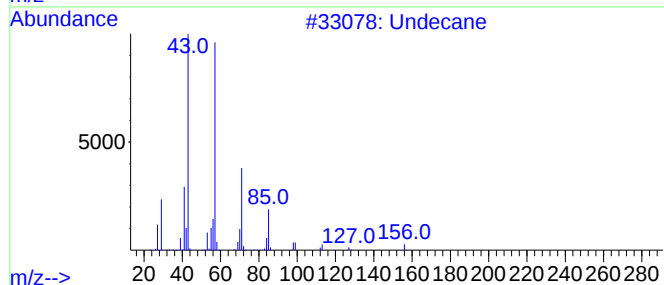
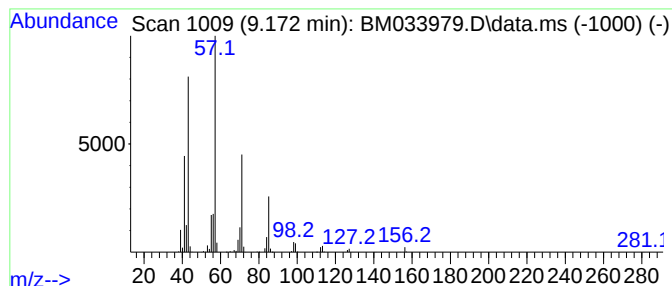
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.172	34.39 ng/ul	2112660	1,4-Dichlorobenzene-d4	7.925

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	94
2	Undecane		156	C11H24	001120-21-4	91
3	Hexadecane		226	C16H34	000544-76-3	86
4	Tridecane		184	C13H28	000629-50-5	86
5	Hexadecane		226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

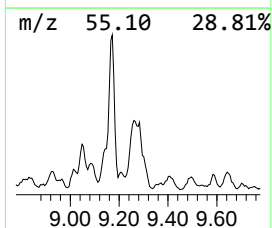
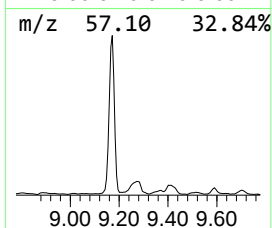
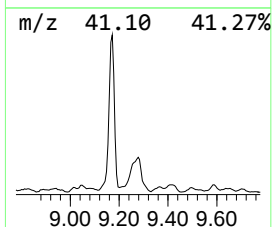
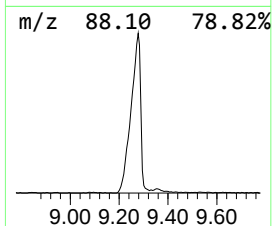
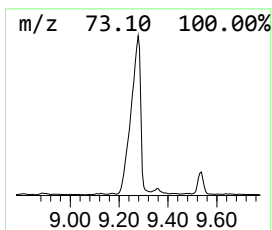
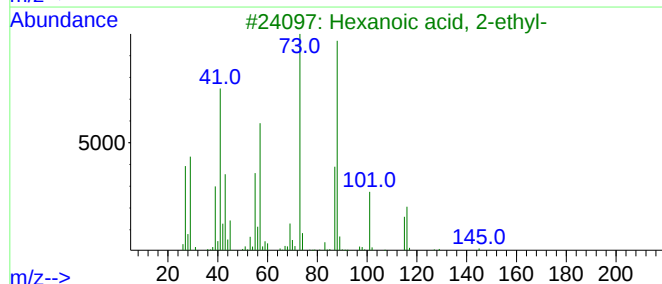
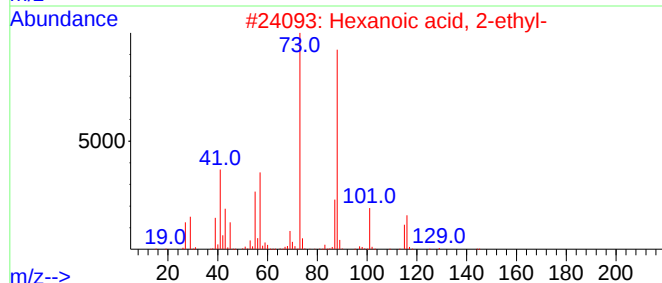
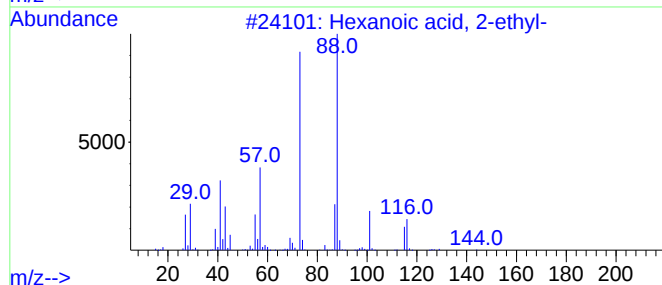
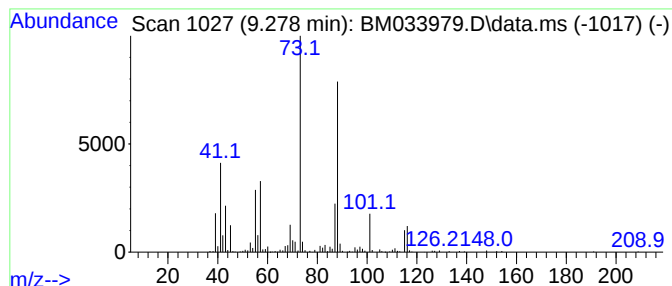
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Hexanoic acid, 2-ethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.278	20.09 ng/ul	1234070	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
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	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

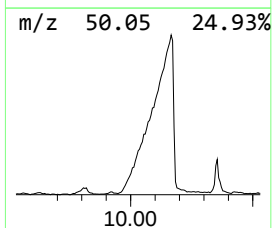
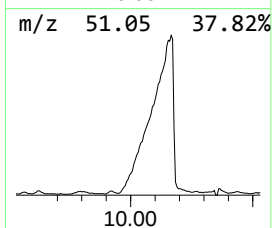
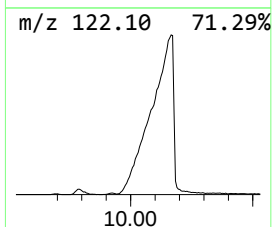
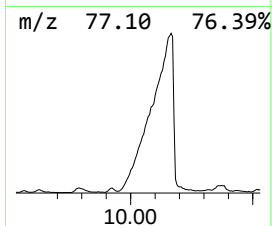
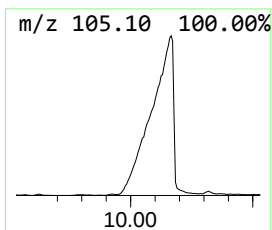
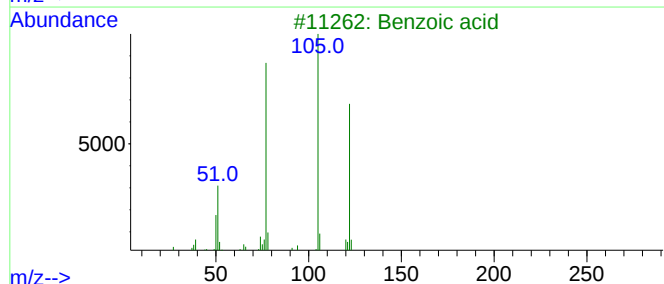
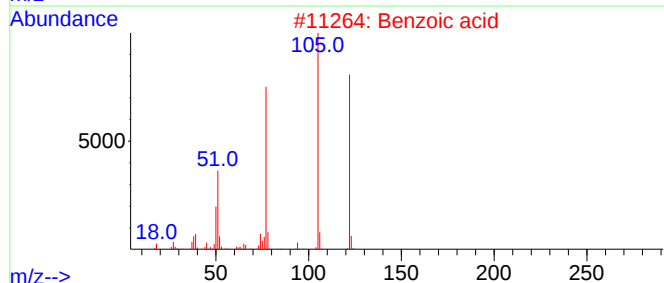
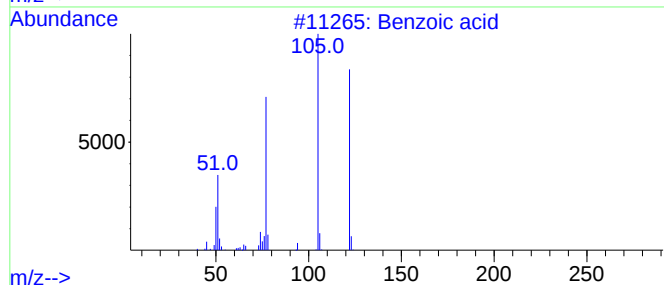
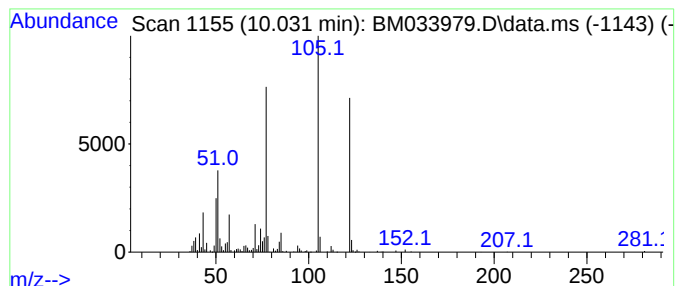
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Benzoic acid Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.031	7.76 ng/ul	717695	Naphthalene-d8	10.736

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	96
2			Benzoic acid	122	C7H6O2	000065-85-0	95
3			Benzoic acid	122	C7H6O2	000065-85-0	91
4			Benzoic acid	122	C7H6O2	000065-85-0	90
5			Cyclobutane-1,1-dicarboxamide, N...	382	C20H18N2O6	1000253-25-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

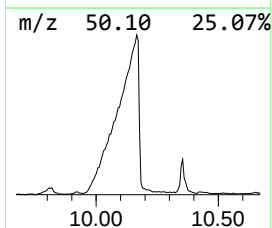
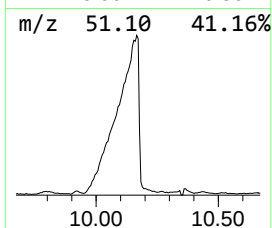
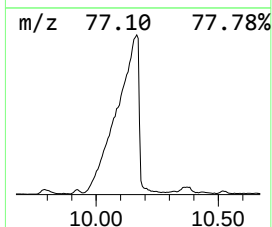
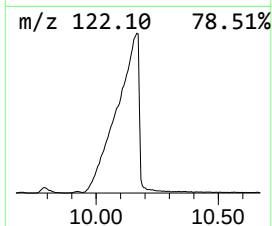
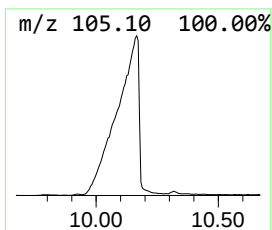
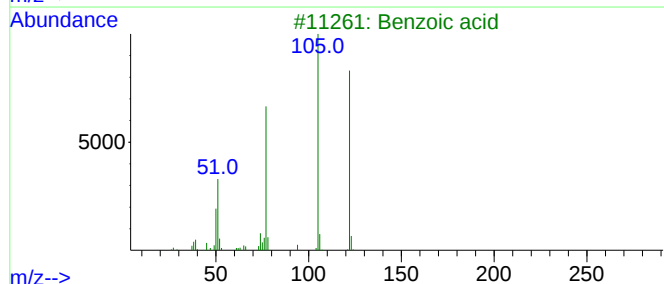
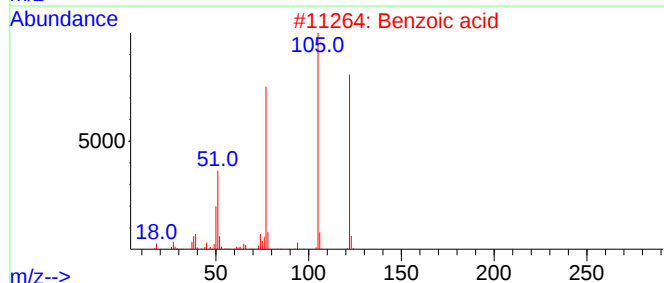
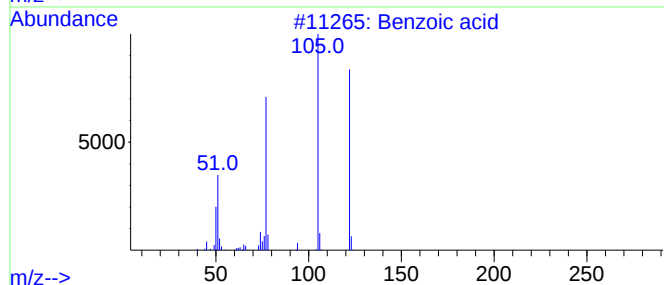
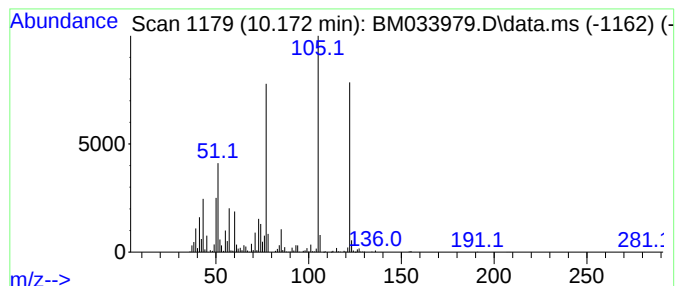
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 Heptanediamide, N,N'-di-ben... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.172	51.17 ng/ul	4734680	Naphthalene-d8	10.736

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	95
2			Benzoic acid	122	C7H6O2	000065-85-0	94
3			Benzoic acid	122	C7H6O2	000065-85-0	93
4			Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	86
5			Cyclobutane-1,1-dicarboxamide, N...	382	C20H18N2O6	1000253-25-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

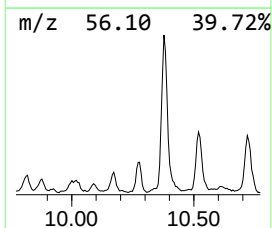
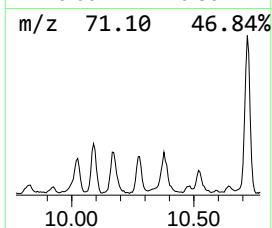
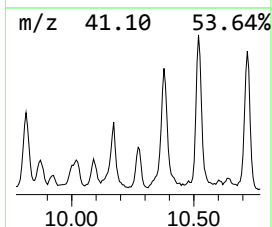
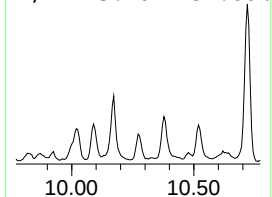
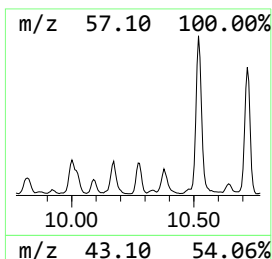
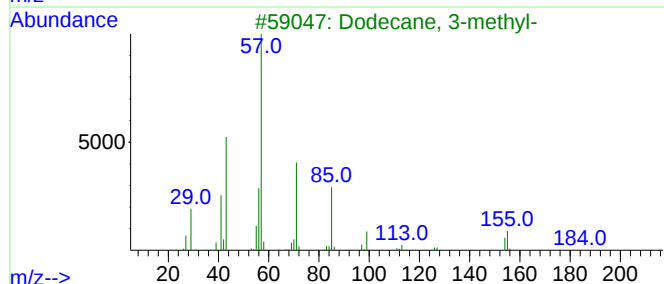
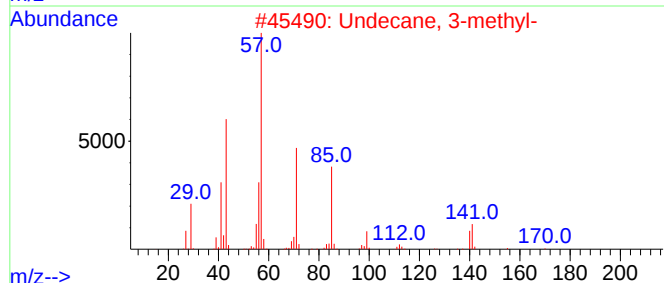
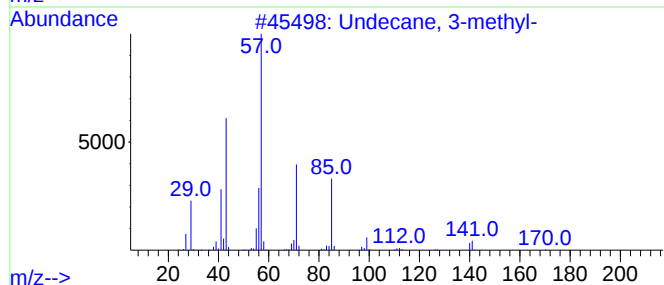
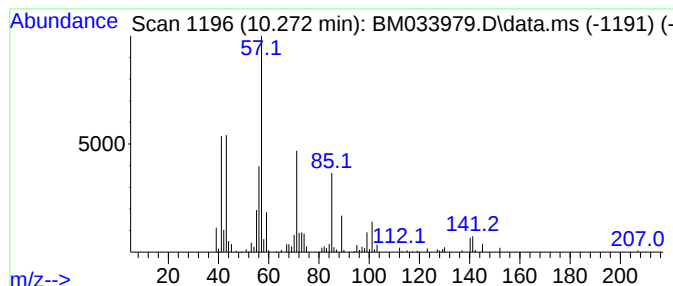
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.272	2.13 ng/ul	196879	Naphthalene-d8	10.736	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3-methyl-	170	C12H26	001002-43-3	59
2	Undecane, 3-methyl-	170	C12H26	001002-43-3	59
3	Dodecane, 3-methyl-	184	C13H28	017312-57-1	59
4	Decane, 3-methyl-	156	C11H24	013151-34-3	50
5	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
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Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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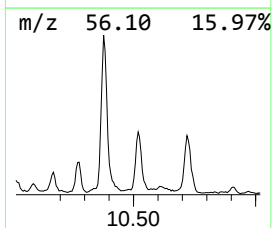
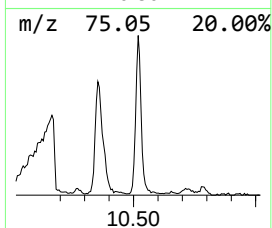
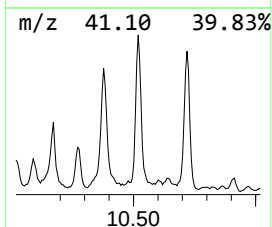
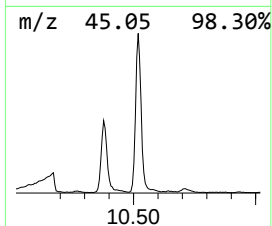
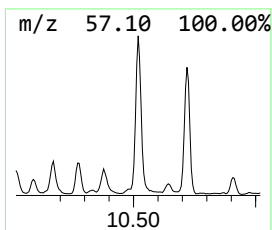
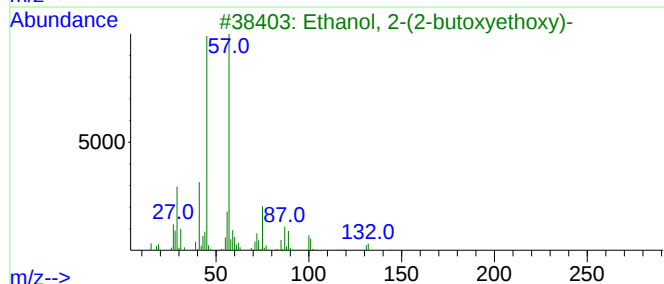
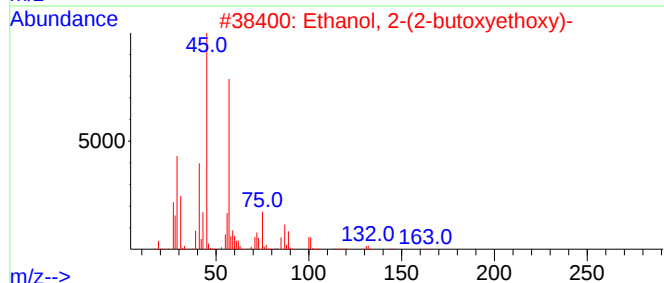
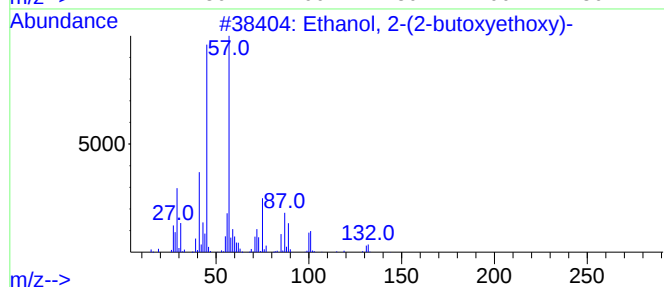
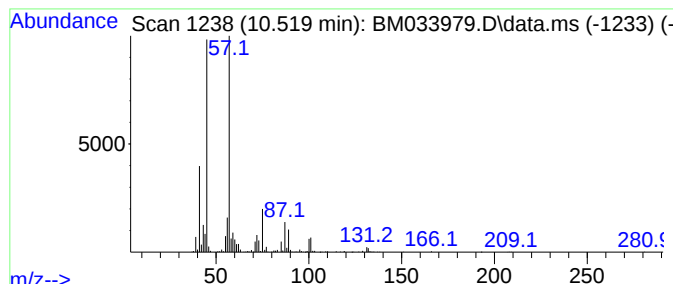
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.519	8.53 ng/ul	789036	Naphthalene-d8	10.736

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

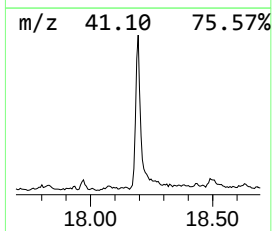
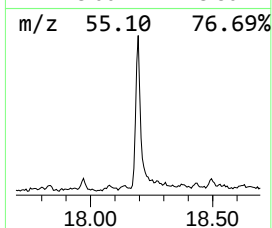
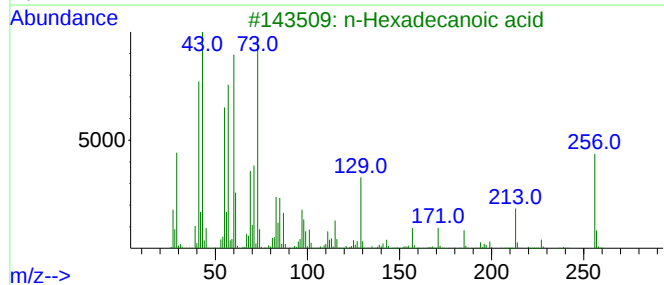
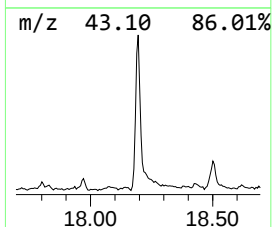
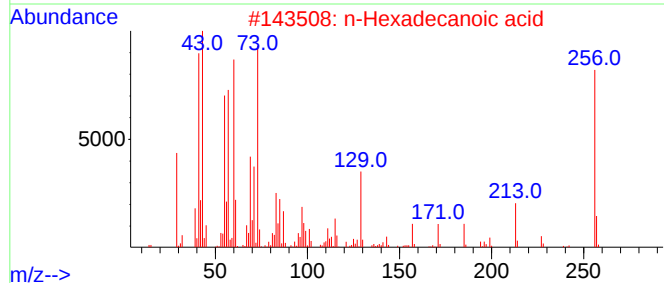
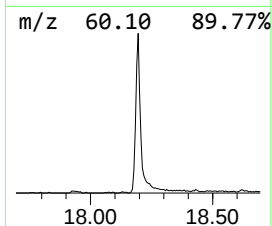
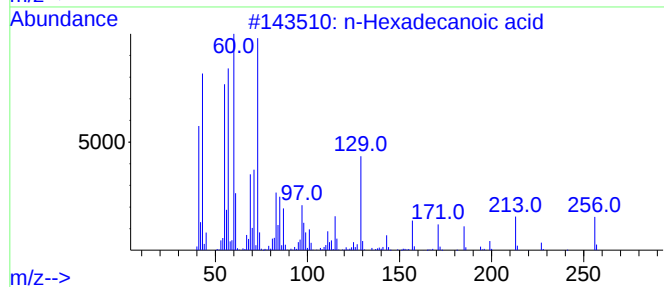
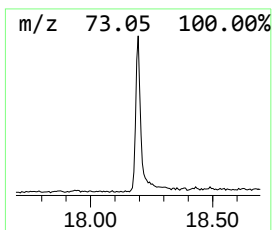
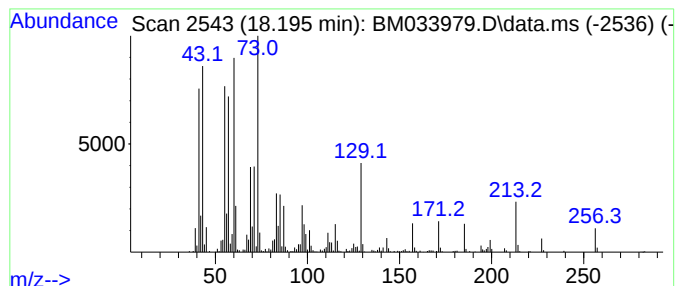
Instrument :
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ClientSampleId :
COVE3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 n-Hexadecanoic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.195	9.19 ng/ul	792460	Phenanthrene-d10	17.295	
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	98
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

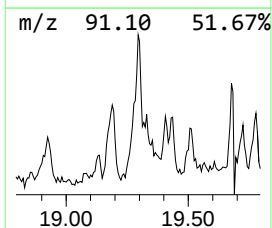
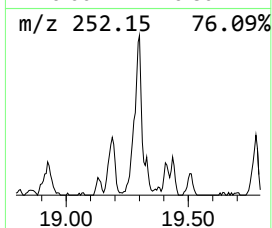
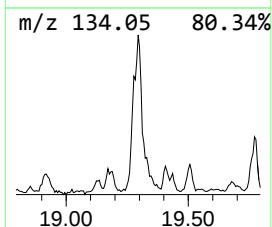
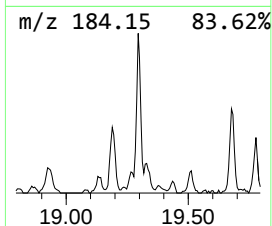
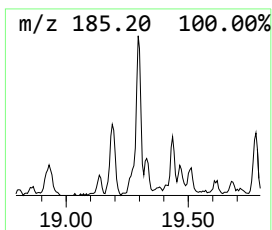
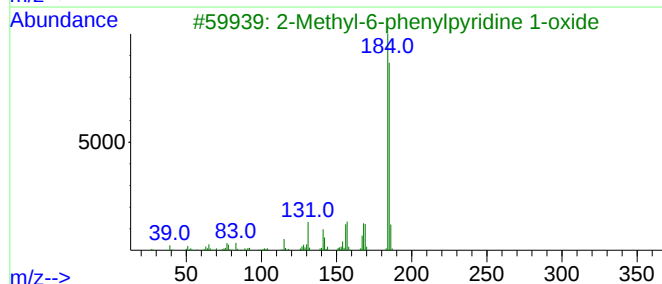
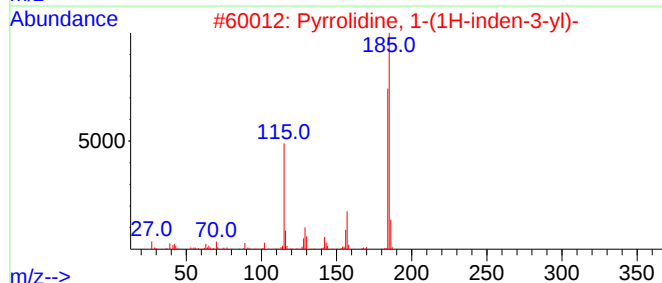
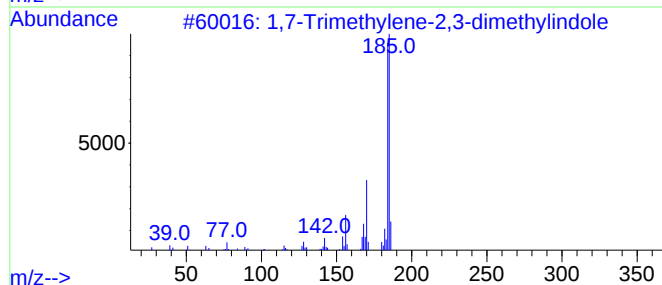
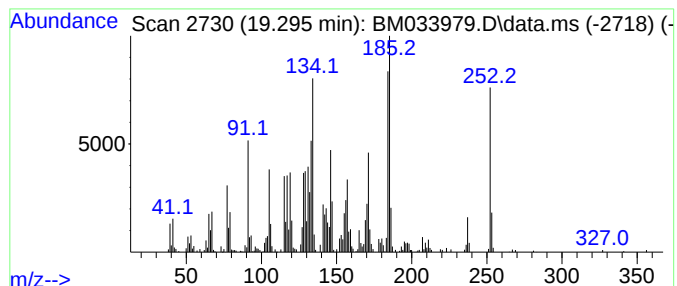
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 unknown-02 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.295	6.75 ng/ul	582052	Phenanthrene-d10	17.295

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,7-Trimethylene-2,3-dimethylindole	185	C13H15N	005825-43-4	38
2			Pyrrolidine, 1-(1H-inden-3-yl)-	185	C13H15N	031554-37-7	30
3			2-Methyl-6-phenylpyridine 1-oxide	185	C12H11NO	020531-89-9	30
4			1,7-Trimethylene-2,3-dimethylindole	185	C13H15N	005825-43-4	27
5			N-Benzylpyrimidin-2-amine	185	C11H11N3	004214-59-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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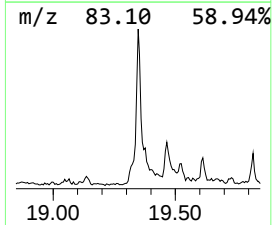
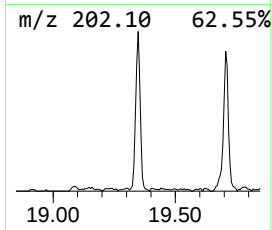
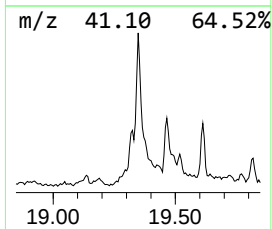
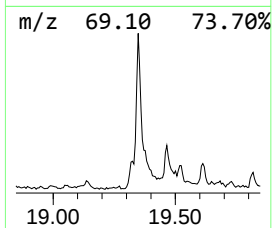
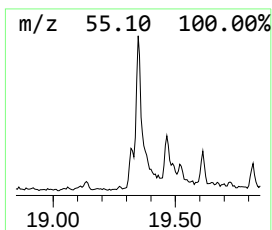
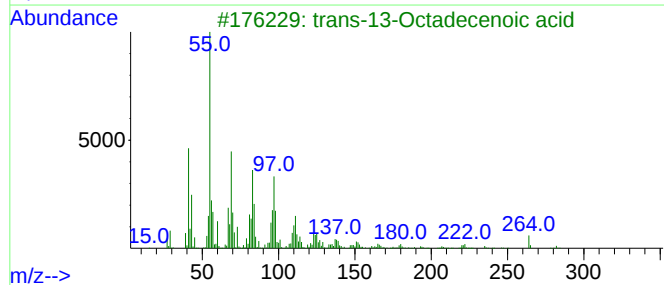
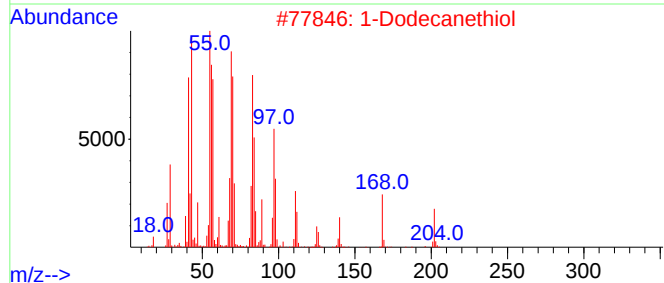
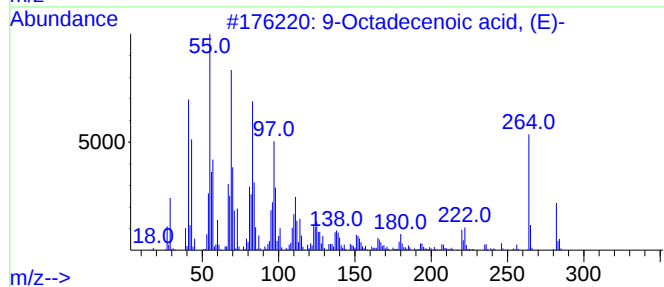
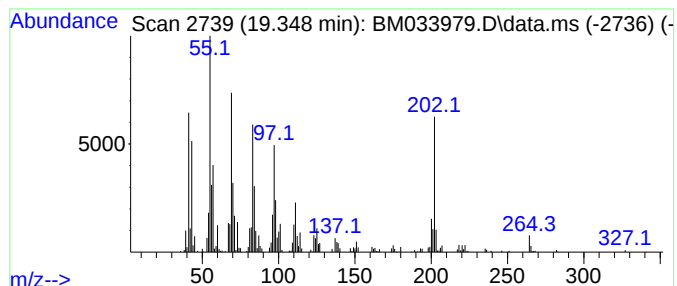
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 9-Octadecenoic acid, (E)- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.348	7.07 ng/ul	609827	Phenanthrene-d10	17.295

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	98
2			1-Dodecanethiol	202	C12H26S	000112-55-0	83
3			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	55
4			3-Bromobenzoic acid, hexadecyl e...	424	C23H37BrO2	070152-90-8	49
5			Acetic acid, trifluoro-, undecyl...	268	C13H23F3O2	053800-01-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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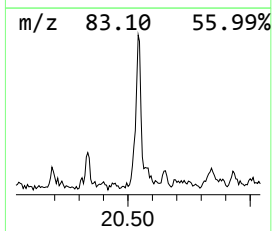
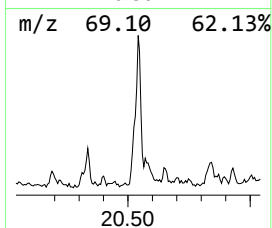
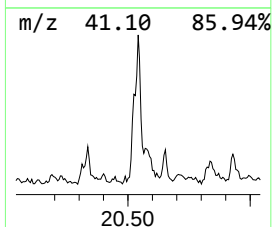
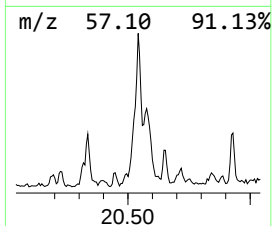
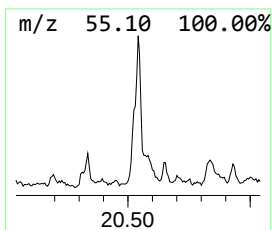
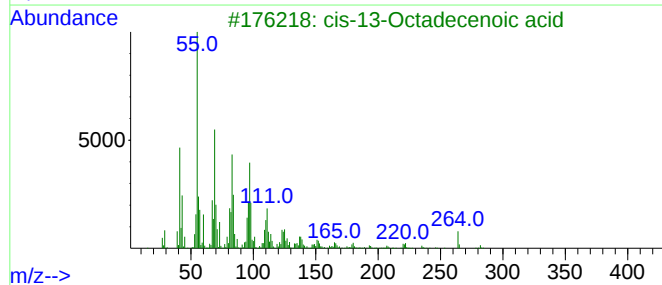
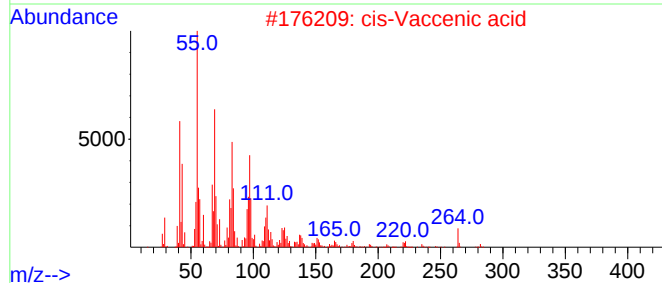
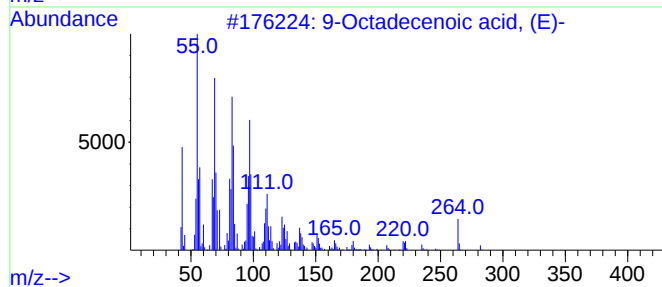
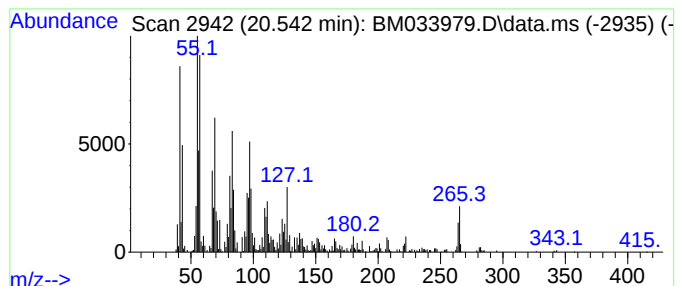
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Quant Title : SVOA CALIBRATION

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

Peak Number 51 cis-Vaccenic acid Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.542	6.16 ng/ul	574018	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	98
2			cis-Vaccenic acid	282	C18H34O2	000506-17-2	93
3			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	92
4			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	91
5			n-Propyl 9-octadecenoate	324	C21H40O2	1000336-71-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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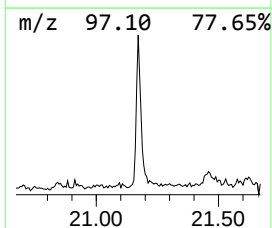
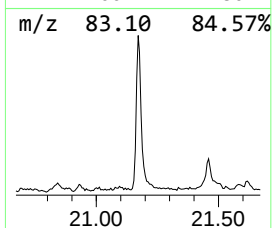
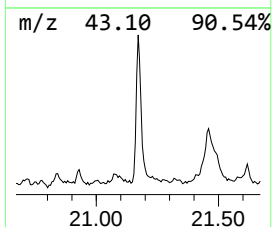
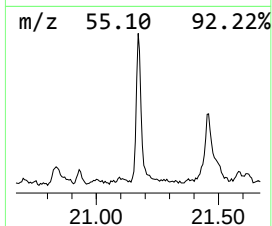
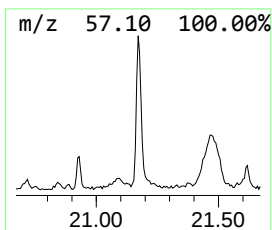
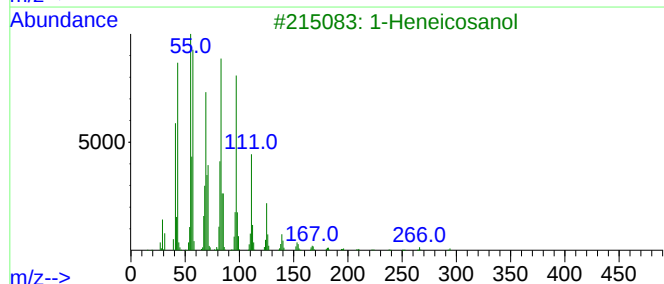
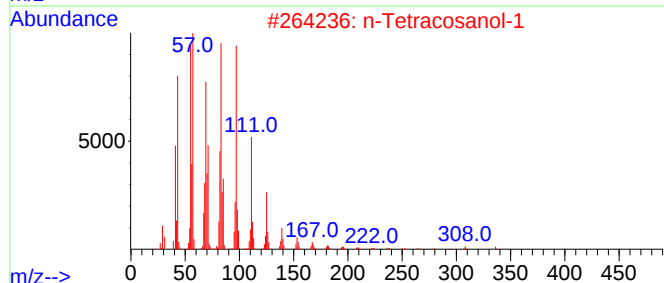
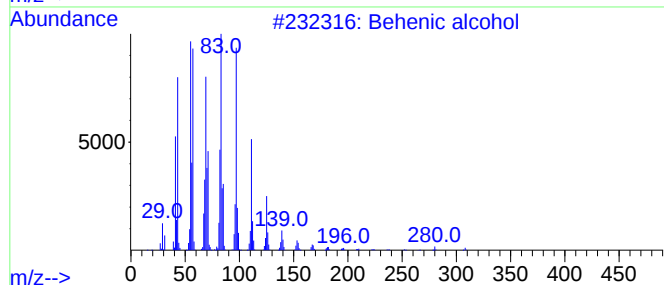
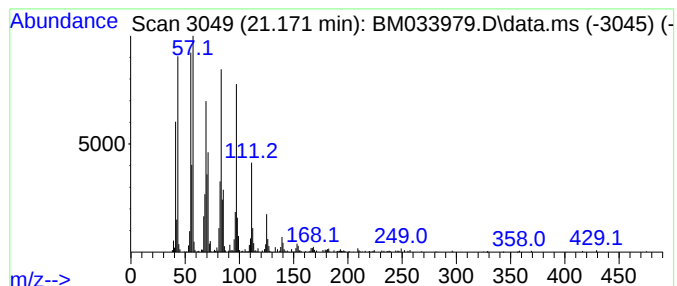
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Quant Title : SVOA CALIBRATION

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

Peak Number 54 Behenic alcohol Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.171	5.22 ng/ul	486207	Chrysene-d12	21.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	95
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	95
3			1-Heneicosanol	312	C21H44O	015594-90-8	95
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

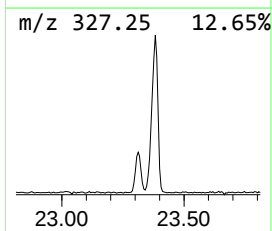
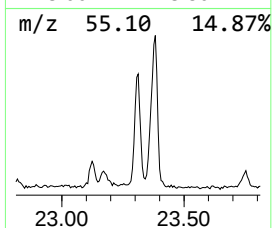
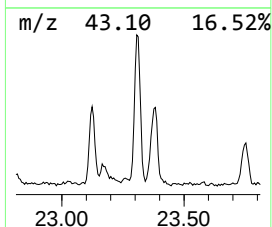
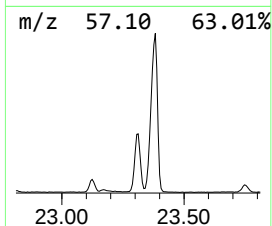
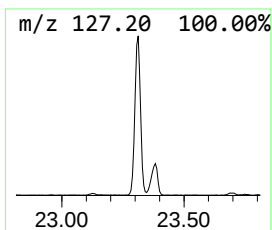
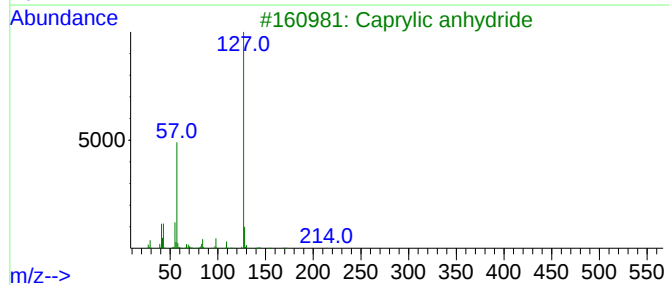
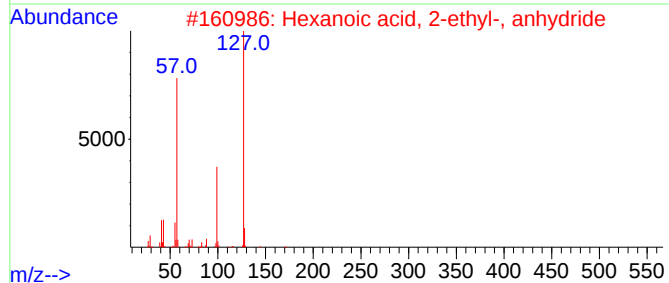
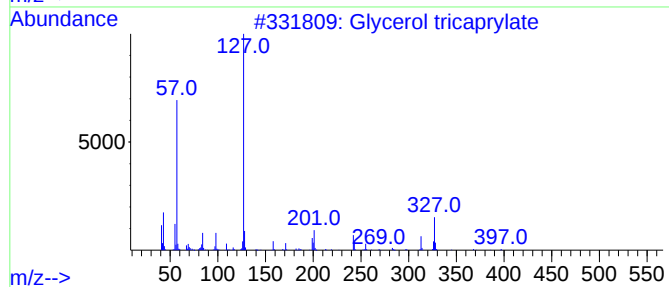
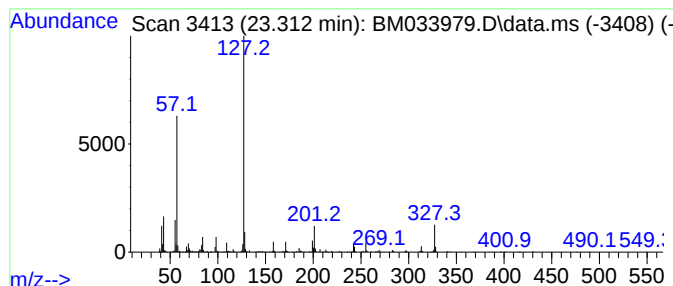
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 Glycerol tricaprylate Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.312	11.70 ng/ul	843319	Perylene-d12	23.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Glycerol tricaprylate	470	C27H50O6	000538-23-8	87
2			Hexanoic acid, 2-ethyl-, anhydride	270	C16H30O3	036765-89-6	50
3			Caprylic anhydride	270	C16H30O3	000623-66-5	50
4			1,2-Dioctanoin	344	C19H36O5	001069-87-0	45
5			Octanoic acid, 1-(hydroxymethyl)...	344	C19H36O5	060514-48-9	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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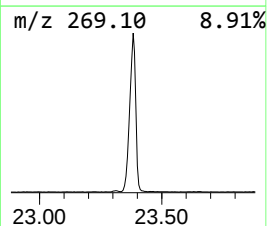
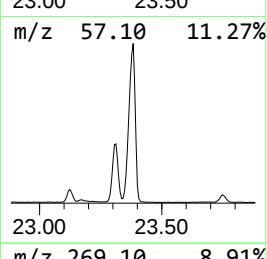
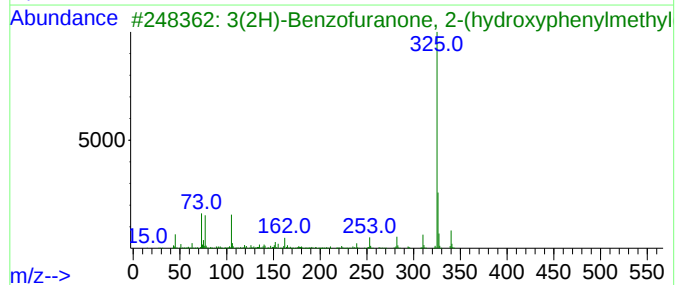
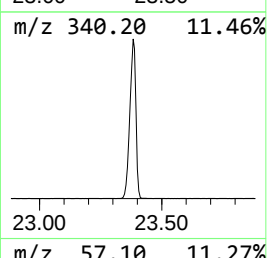
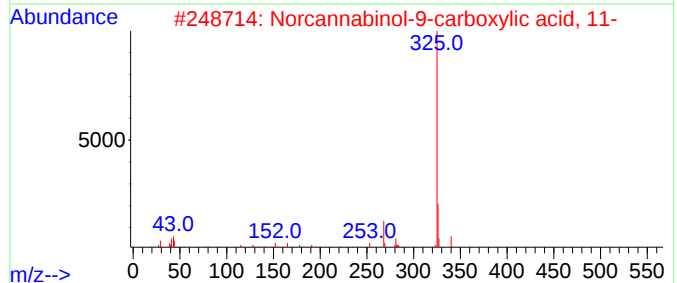
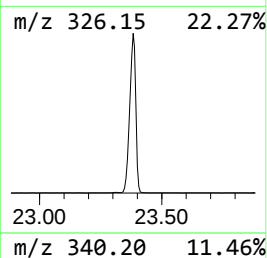
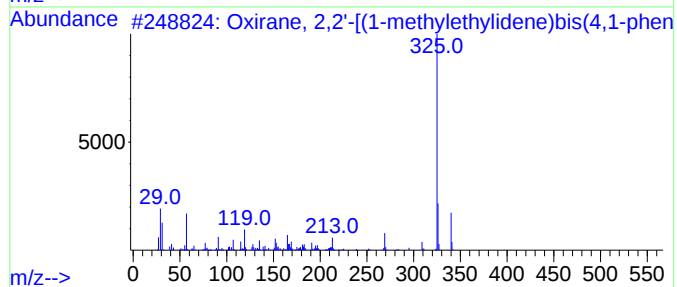
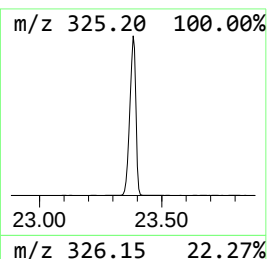
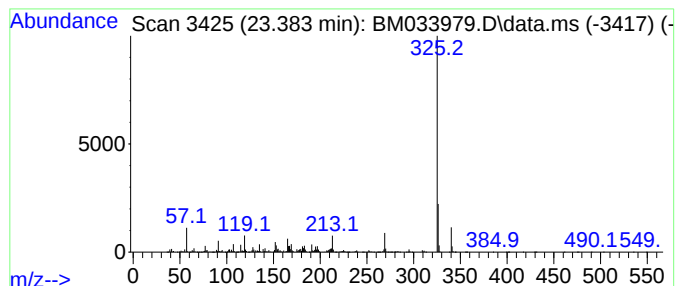
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 Oxirane, 2,2'-[(1-methyleth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.383	177.62 ng/ul	12806000	Perylene-d12	23.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, 2,2'-[(1-methylethylidene)bis(4,1-phen	340	C21H24O4	001675-54-3	95
2			Norcannabinol-9-carboxylic acid, 11-	340	C21H24O4	053989-32-5	80
3			3(2H)-Benzofuranone, 2-(hydroxyphenylmethyl	340	C19H20O4Si	1000503-53-4	72
4			Methyl 4,6-dihydroxy-2,3-dimethyl-5-(2-methyl-2-propenyl)-2H-pyran-2-carboxylate	340	C16H28O4Si2	1000502-58-9	72
5			2,6-Bis(1,1-dimethylethyl)cyclohexanone	325	C21H27NO2	017119-01-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

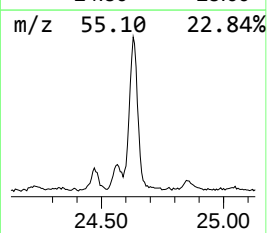
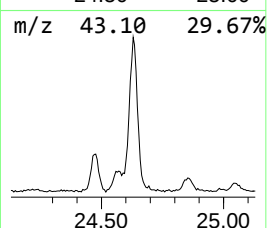
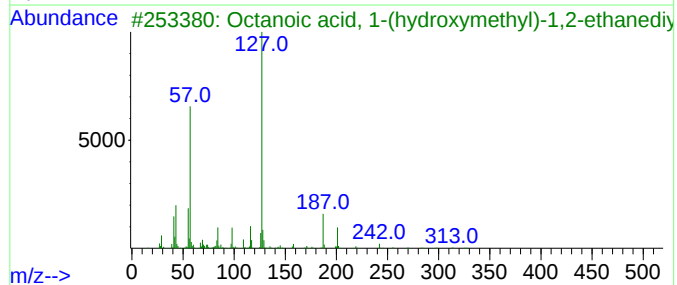
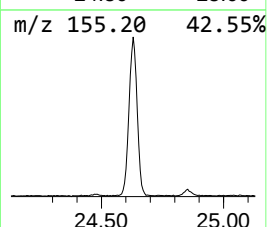
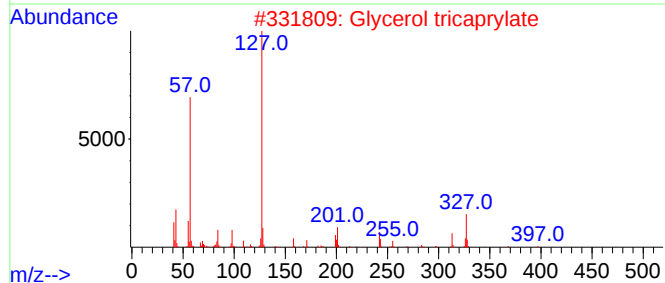
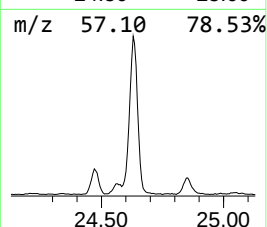
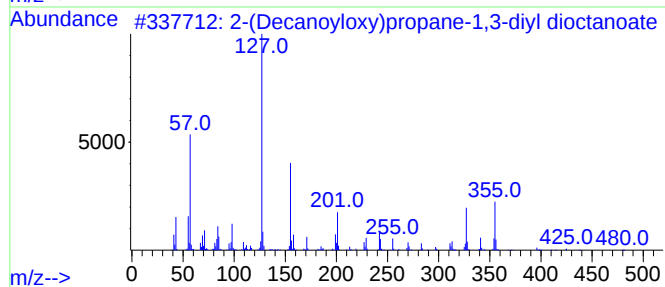
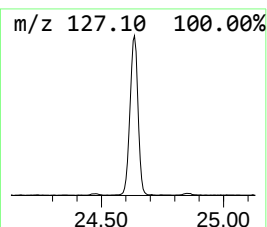
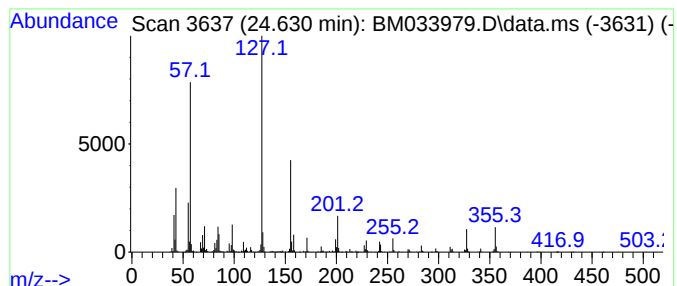
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 2-(Decanoyloxy)propane-1,3-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.630	24.37 ng/ul	1757210	Perylene-d12	23.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(Decanoyloxy)propane-1,3-diyl ...	498	C29H54O6	033368-87-5	72		
2	Glycerol tricaprylate	470	C27H50O6	000538-23-8	50		
3	Octanoic acid, 1-(hydroxymethyl)...	344	C19H36O5	060514-48-9	50		
4	1-(Nonanoyloxy)-3-(octanoyloxy)p...	512	C30H56O6	1000413-78-7	50		
5	3-Fluoro-5-methoxypyridine	127	C6H6FNO	1060801-62-8	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

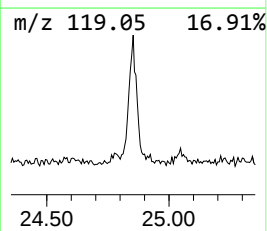
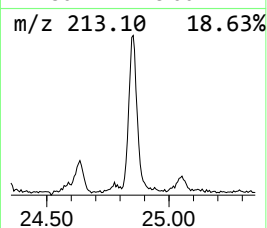
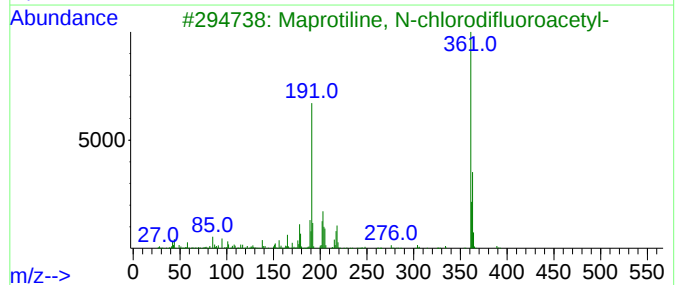
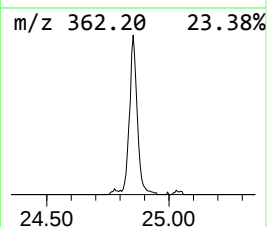
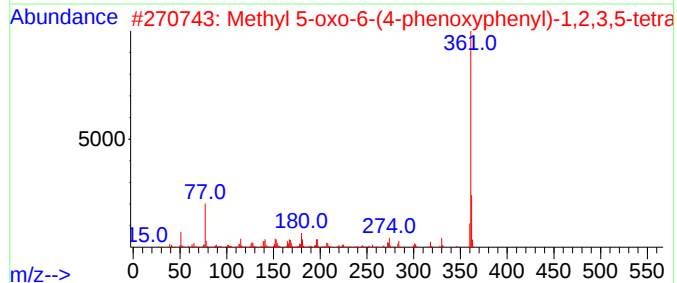
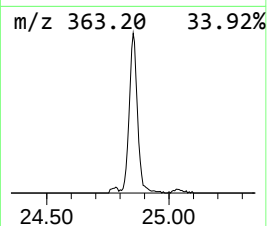
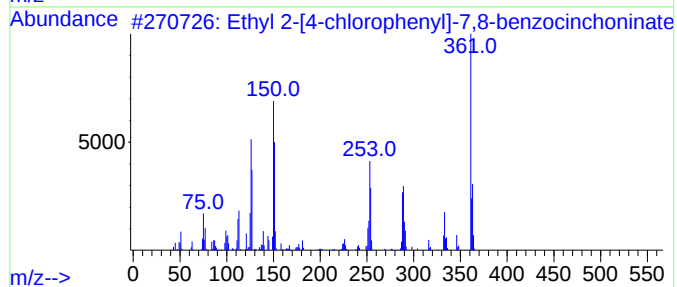
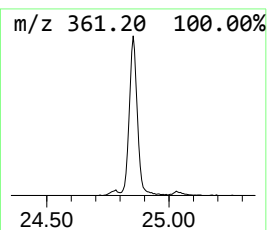
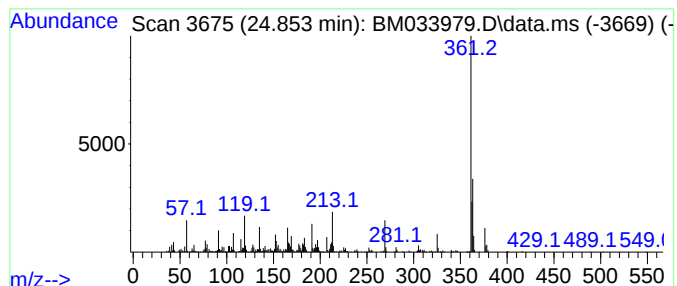
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 Ethyl 2-[4-chlorophenyl]-7,... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.853	13.78 ng/ul	993337	Perylene-d12	23.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl 2-[4-chlorophenyl]-7,8-ben...	361	C22H16ClNO2	1000256-80-8	94
2			Methyl 5-oxo-6-(4-phenoxyphenyl)...	361	C22H19NO4	1000482-02-4	50
3			Maprotiline, N-chlorodifluoroace...	389	C22H22ClF2NO	1000448-25-9	45
4			Acifluorfen	361	C14H7ClF3NO5	050594-66-6	45
5			3-Bromo-3',5,5'-trihydroxy-2,2'...	440	C20H9BrO7	088381-89-9	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033979.D
Acq On : 02 Feb 2022 18:26
Operator : CG/JU
Sample : N1355-02
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COVE3

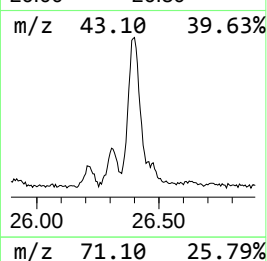
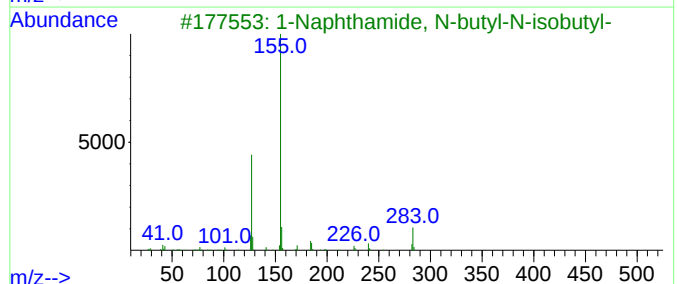
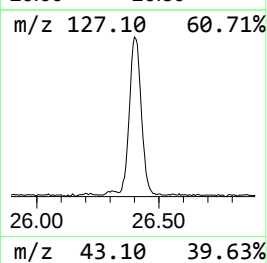
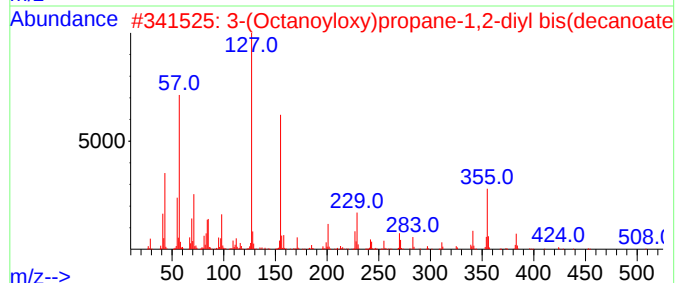
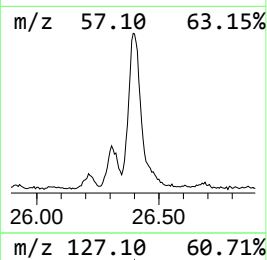
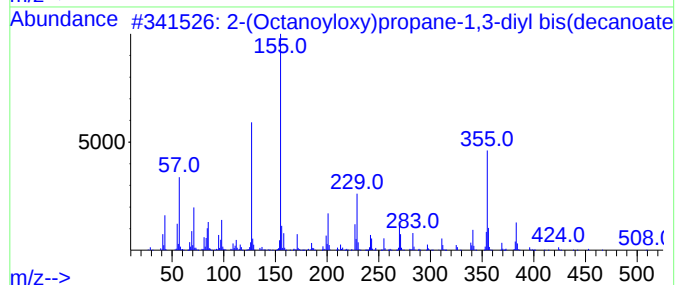
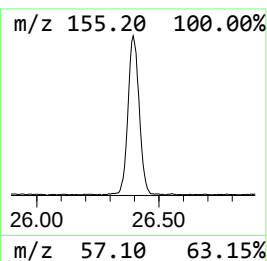
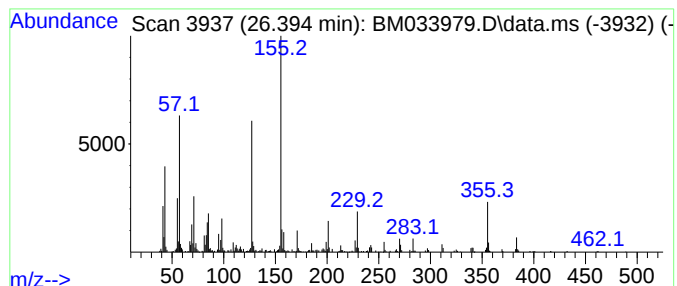
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 2-(Octanoyloxy)propane-1,3-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.394	15.15 ng/ul	1092340	Perylene-d12	23.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(Octanoyloxy)propane-1,3-diyl ...	526	C31H58O6	033368-86-4	90		
2	3-(Octanoyloxy)propane-1,2-diyl ...	526	C31H58O6	082426-88-8	53		
3	1-Naphthamide, N-butyl-N-isobutyl-	283	C19H25NO	1000415-71-5	49		
4	1-(Nonanoyloxy)-3-(octanoyloxy)p...	512	C30H56O6	1000413-78-7	49		
5	Phenylpropionic acid, .alpha.-am...	229	C10H12FN04	1000126-07-3	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033979.D
 Acq On : 02 Feb 2022 18:26
 Operator : CG/JU
 Sample : N1355-02
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COVE3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.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
Methyl Isobutyl...	3.649	45.6	ng/ul	2802050	1	7.925	1228750	20.0
2-Pentanol, 4-m...	3.878	10.6	ng/ul	651273	1	7.925	1228750	20.0
Butanoic acid	4.060	22.1	ng/ul	1357340	1	7.925	1228750	20.0
Ethanol, 2-prop...	4.472	25.9	ng/ul	1593050	1	7.925	1228750	20.0
unknown-01	4.890	5.2	ng/ul	316840	1	7.925	1228750	20.0
2-Pentanone, 4-...	5.049	342.7	ng/ul	21054000	1	7.925	1228750	20.0
Ethylbenzene	5.407	12.1	ng/ul	742136	1	7.925	1228750	20.0
Benzene, 1,3-di...	5.543	21.1	ng/ul	1293650	1	7.925	1228750	20.0
2-Heptanone	5.754	7.4	ng/ul	454530	1	7.925	1228750	20.0
o-Xylene	5.925	17.2	ng/ul	1057160	1	7.925	1228750	20.0
Ethanol, 2-butoxy-	6.001	48.9	ng/ul	3003610	1	7.925	1228750	20.0
(DEL) Alkane: C...	6.190	2.4	ng/ul	145094	1	7.925	1228750	20.0
(DEL) Alkane: S...	6.478	3.9	ng/ul	238971	1	7.925	1228750	20.0
(DEL) Alkane: C...	6.560	3.0	ng/ul	183947	1	7.925	1228750	20.0
(DEL) Alkane: S...	6.884	2.1	ng/ul	132077	1	7.925	1228750	20.0
(DEL) Alkane: S...	6.919	4.2	ng/ul	257114	1	7.925	1228750	20.0
(DEL) Alkane: S...	6.978	6.9	ng/ul	425255	1	7.925	1228750	20.0
Benzene, 1-ethy...	7.013	7.7	ng/ul	470732	1	7.925	1228750	20.0
2-Propanol, 1-(...	7.531	54.0	ng/ul	3316490	1	7.925	1228750	20.0
2-Propanol, 1-(...	7.737	22.3	ng/ul	1368800	1	7.925	1228750	20.0
Mesitylene	8.048	4.8	ng/ul	292794	1	7.925	1228750	20.0
Benzyl alcohol	8.213	484.5	ng/ul	29767200	1	7.925	1228750	20.0
(DEL) Alkane: S...	8.466	9.3	ng/ul	571173	1	7.925	1228750	20.0
(DEL) Alkane: S...	8.525	4.2	ng/ul	256334	1	7.925	1228750	20.0
(DEL) Alkane: S...	9.172	34.4	ng/ul	2112660	1	7.925	1228750	20.0
Hexanoic acid, ...	9.278	20.1	ng/ul	1234070	1	7.925	1228750	20.0
Benzoic acid	10.031	7.8	ng/ul	717695	2	10.736	1850590	20.0
Heptanediamide,...	10.172	51.2	ng/ul	4734680	2	10.736	1850590	20.0
(DEL) Alkane: S...	10.272	2.1	ng/ul	196879	2	10.736	1850590	20.0
Ethanol, 2-(2-b...	10.519	8.5	ng/ul	789036	2	10.736	1850590	20.0
n-Hexadecanoic ...	18.195	9.2	ng/ul	792460	4	17.295	1724600	20.0
unknown-02	19.295	6.8	ng/ul	582052	4	17.295	1724600	20.0
9-Octadecenoic ...	19.348	7.1	ng/ul	609827	4	17.295	1724600	20.0
cis-Vaccenic acid	20.542	6.2	ng/ul	574018	5	21.459	1863690	20.0
Behenic alcohol	21.171	5.2	ng/ul	486207	5	21.459	1863690	20.0
Glycerol tricap...	23.312	11.7	ng/ul	843319	6	23.853	1441950	20.0
Oxirane, 2,2'-[...	23.383	177.6	ng/ul	12806000	6	23.853	1441950	20.0
2-(Decanoyloxy)...	24.630	24.4	ng/ul	1757210	6	23.853	1441950	20.0
Ethyl 2-[4-chlo...	24.853	13.8	ng/ul	993337	6	23.853	1441950	20.0
2-(Octanoyloxy)...	26.394	15.2	ng/ul	1092340	6	23.853	1441950	20.0