

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
 Data File : BP016331.D
 Acq On : 19 Jul 2023 19:20
 Operator : MA/JU
 Sample : 03472-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A46Q5

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_P\Methods\SFAM-EPA-BP071223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016331.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.276	12	15	21	rVB2	15062	21126	0.81%	0.074%
2	3.746	93	95	110	rBV	67439	186513	7.12%	0.654%
3	4.040	141	145	152	rVB	30427	49568	1.89%	0.174%
4	4.217	172	175	184	rBV	41631	120709	4.61%	0.423%
5	4.581	232	237	242	rBV8	7958	13732	0.52%	0.048%
6	4.799	272	274	277	rBV4	2588	2781	0.11%	0.010%
7	5.358	363	369	375	rBV8	12917	25427	0.97%	0.089%
8	5.817	437	447	456	rBV2	214548	482207	18.41%	1.690%
9	5.899	456	461	477	rVB2	48759	159418	6.09%	0.559%
10	6.105	492	496	499	rVB5	2504	3499	0.13%	0.012%
11	6.181	504	509	516	rVV2	44859	76425	2.92%	0.268%
12	6.293	526	528	532	rVB5	2005	2773	0.11%	0.010%
13	6.581	570	577	597	rBV	266372	706949	26.99%	2.478%
14	7.175	671	678	693	rBV2	101353	413738	15.80%	1.450%
15	7.293	693	698	718	rVB	127199	363708	13.89%	1.275%
16	7.499	727	733	759	rVB	170354	450244	17.19%	1.578%
17	7.952	802	810	825	rBV	217718	608582	23.24%	2.133%
18	8.058	825	828	839	rVV3	33419	89452	3.42%	0.313%
19	8.164	841	846	850	rVB6	6969	14156	0.54%	0.050%
20	8.263	856	863	880	rBV	314984	658514	25.14%	2.308%
21	8.493	900	902	905	rVB4	2723	2741	0.10%	0.010%
22	8.687	933	935	939	rVB5	3516	3121	0.12%	0.011%
23	8.793	945	953	957	rBV7	35663	108109	4.13%	0.379%
24	9.081	1000	1002	1009	rVB8	2513	3488	0.13%	0.012%
25	9.187	1013	1020	1049	rBV	136071	523569	19.99%	1.835%
26	9.458	1063	1066	1075	rVB10	5492	9950	0.38%	0.035%
27	9.634	1092	1096	1098	rVB5	2087	3066	0.12%	0.011%
28	9.916	1136	1144	1166	rBV3	83924	437977	16.72%	1.535%
29	10.475	1230	1239	1240	rBV8	7257	13819	0.53%	0.048%
30	10.528	1241	1248	1275	rVV3	87674	499619	19.08%	1.751%
31	10.781	1284	1291	1321	rVB	332015	989333	37.78%	3.467%
32	11.087	1334	1343	1355	rBV3	20533	94805	3.62%	0.332%
33	11.269	1371	1374	1375	rBV3	2912	3200	0.12%	0.011%
34	11.675	1440	1443	1448	rVB7	2481	2897	0.11%	0.010%
35	11.928	1481	1486	1488	rBV6	2813	3694	0.14%	0.013%
36	12.063	1506	1509	1510	rBV3	3427	4003	0.15%	0.014%

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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_P\Methods\SFAM-EPA-BP071223.MA.M
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37	12.446	1568	1574	1579	rBV9	6194	12035	0.46%	0.042%
38	12.746	1618	1625	1629	rBV9	5691	14417	0.55%	0.051%
39	12.816	1632	1637	1639	rVV	14289	25892	0.99%	0.091%
40	12.846	1639	1642	1650	rVB	14427	28005	1.07%	0.098%

41	13.199	1699	1702	1703	rBV3	3664	3367	0.13%	0.012%
42	13.493	1746	1752	1762	rBV6	10043	23087	0.88%	0.081%
43	14.034	1835	1844	1873	rBV	489118	1265099	48.31%	4.434%
44	14.304	1883	1890	1906	rBV	693735	1318814	50.36%	4.622%
45	14.446	1912	1914	1921	rVB8	8128	11652	0.44%	0.041%

46	14.604	1934	1941	1968	rVB	818369	1462442	55.84%	5.125%
47	15.010	2002	2010	2019	rBV6	54102	252089	9.63%	0.883%
48	15.404	2076	2077	2084	rVB7	5208	4567	0.17%	0.016%
49	15.616	2106	2113	2134	rBV	754157	1808228	69.05%	6.337%
50	15.834	2143	2150	2173	rBV4	90517	451833	17.25%	1.584%

51	15.998	2173	2178	2191	rVB	121253	213649	8.16%	0.749%
52	16.145	2201	2203	2210	rVB7	7114	13345	0.51%	0.047%
53	16.251	2219	2221	2231	rVB7	4781	9870	0.38%	0.035%
54	16.545	2267	2271	2278	rVB9	7752	14639	0.56%	0.051%
55	16.604	2278	2281	2285	rBV6	3282	5351	0.20%	0.019%

56	16.781	2306	2311	2312	rBV5	5316	5617	0.21%	0.020%
57	16.922	2329	2335	2344	rBV5	14159	32912	1.26%	0.115%
58	17.045	2352	2356	2358	rBV5	4521	6101	0.23%	0.021%
59	17.092	2359	2364	2365	rBV5	10653	14141	0.54%	0.050%
60	17.257	2385	2392	2403	rBV5	17416	44518	1.70%	0.156%

61	17.369	2405	2411	2424	rBV	1154878	1895616	72.38%	6.643%
62	17.481	2424	2430	2450	rVV2	671196	1610025	61.48%	5.643%
63	17.610	2450	2452	2456	rVB4	6927	8806	0.34%	0.031%
64	18.010	2515	2520	2523	rBV7	11107	17753	0.68%	0.062%
65	18.087	2531	2533	2535	rBV2	4811	3985	0.15%	0.014%

66	18.228	2552	2557	2568	rBV	273125	502517	19.19%	1.761%
67	18.422	2585	2590	2596	rBV4	27225	76014	2.90%	0.266%
68	18.834	2654	2660	2662	rBV7	12375	25390	0.97%	0.089%
69	18.975	2680	2684	2695	rBV9	29025	96491	3.68%	0.338%
70	19.286	2734	2737	2740	rBV3	17542	20167	0.77%	0.071%

71	19.363	2748	2750	2754	rVB	25334	28789	1.10%	0.101%
72	19.457	2762	2766	2773	rBV	121459	199327	7.61%	0.699%
73	19.710	2803	2809	2823	rBV	1431490	2618878	100.00%	9.178%
74	21.116	3045	3048	3050	rBV2	53568	64058	2.45%	0.224%
75	21.163	3051	3056	3067	rVB4	155607	453224	17.31%	1.588%

76	21.475	3104	3109	3120	rBV2	1184845	2421521	92.46%	8.487%
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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_P\Methods\SFAM-EPA-BP071223.MA.M
Title : SVOA CALIBRATION

77	23.792	3498	3503	3520	rVB2	755162	1883343	71.91%	6.600%
78	23.957	3525	3531	3556	rVB	817490	2413277	92.15%	8.458%

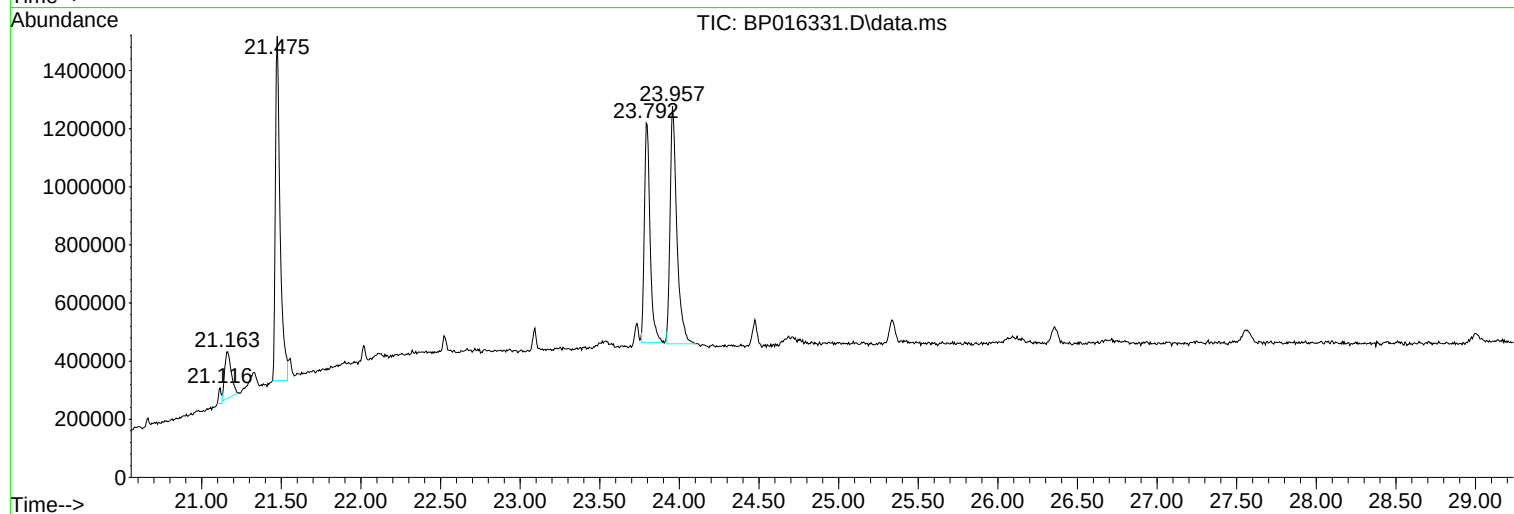
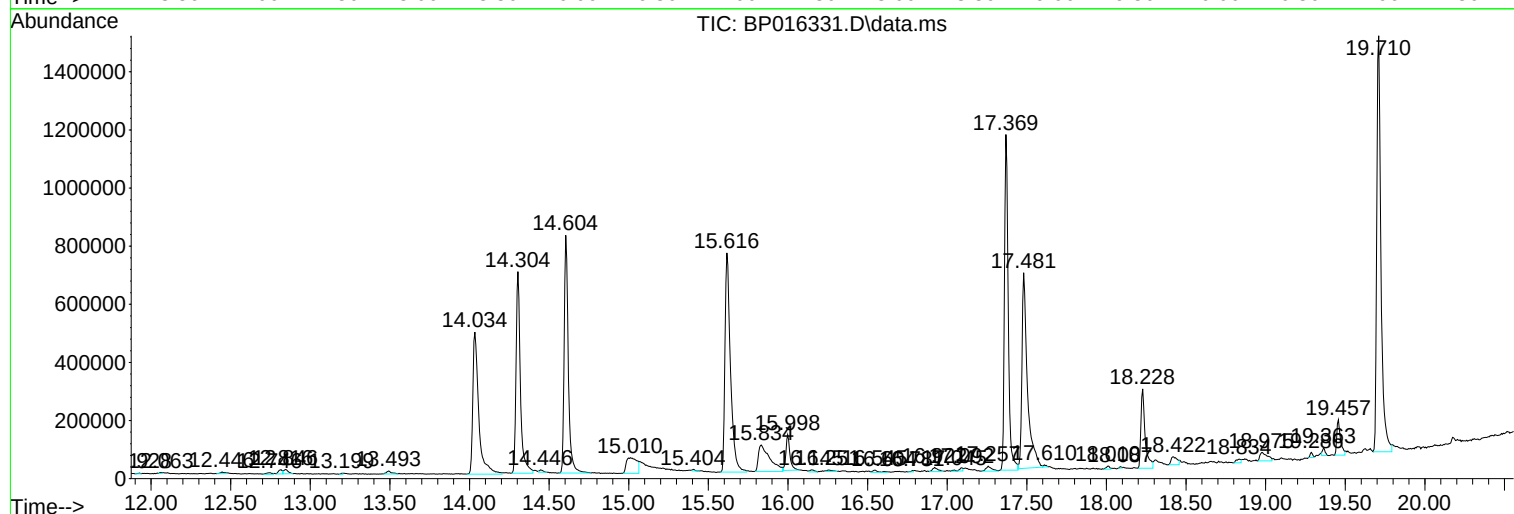
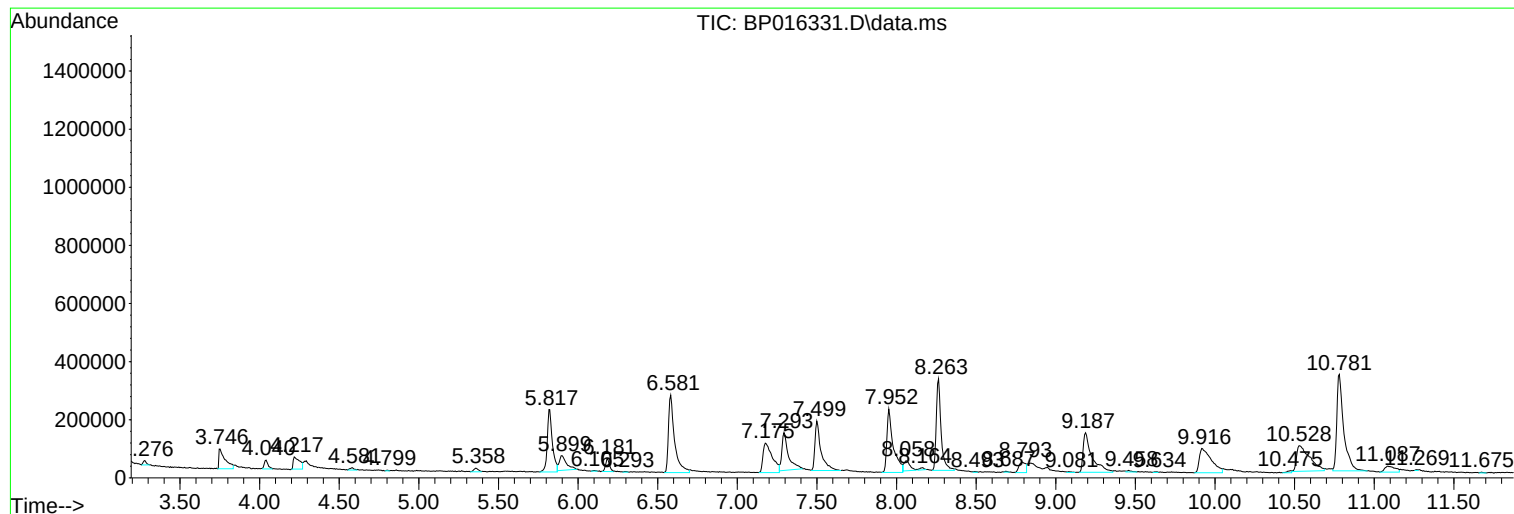
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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



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Sample : 03472-11
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ALS Vial : 14 Sample Multiplier: 1

Instrument :
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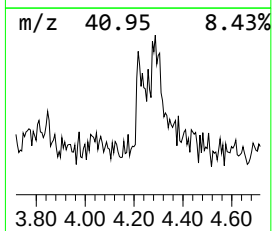
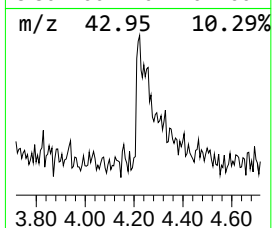
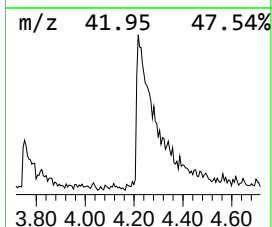
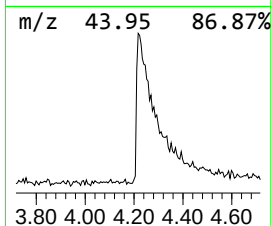
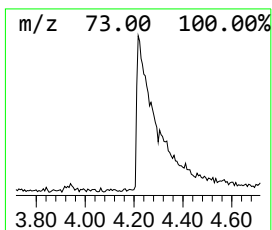
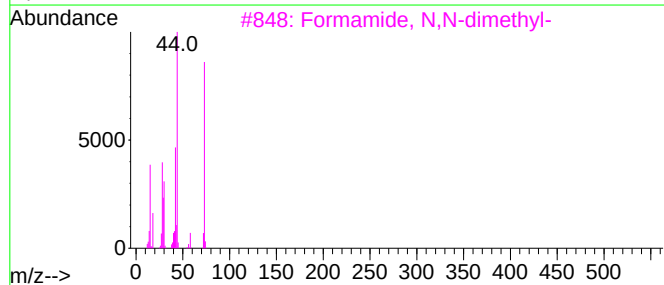
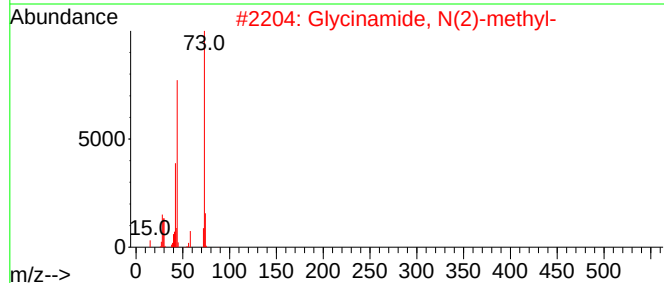
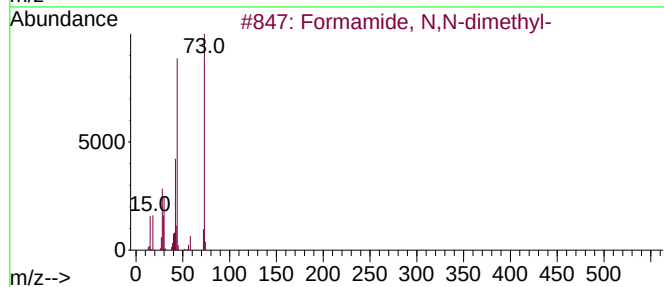
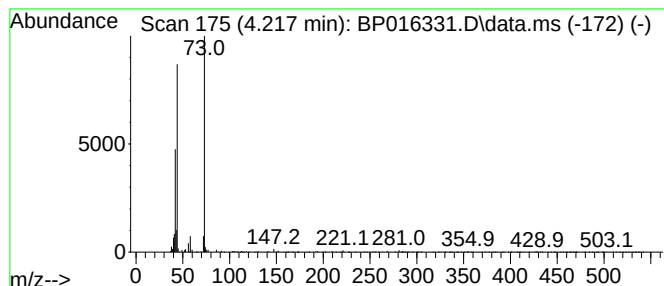
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Formamide, N,N-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.217	3.97 ng/ul	120709	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	72
2			Glycinamide, N(2)-methyl-	88	C3H8N2O	1000452-55-9	56
3			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	56
4			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	38
5			Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	9



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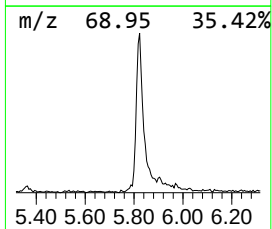
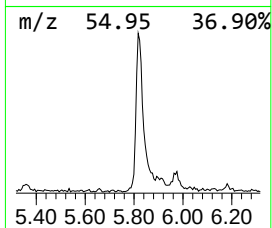
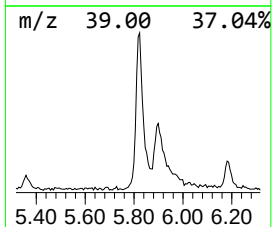
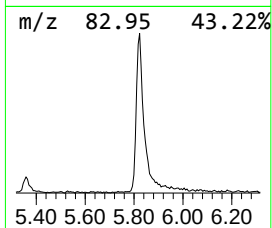
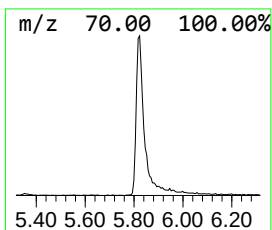
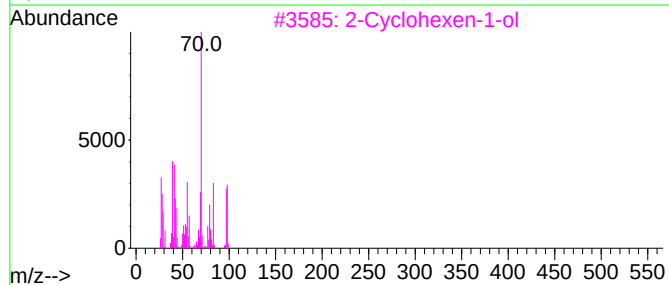
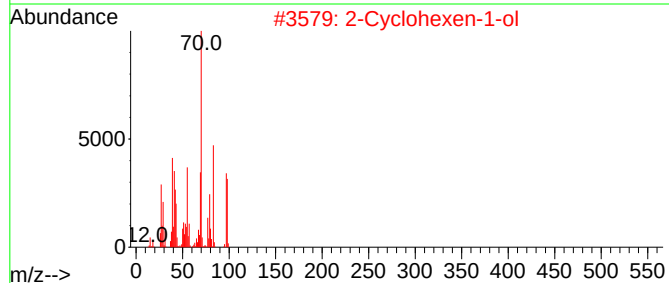
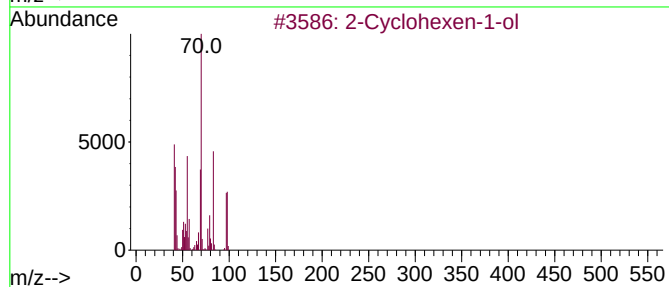
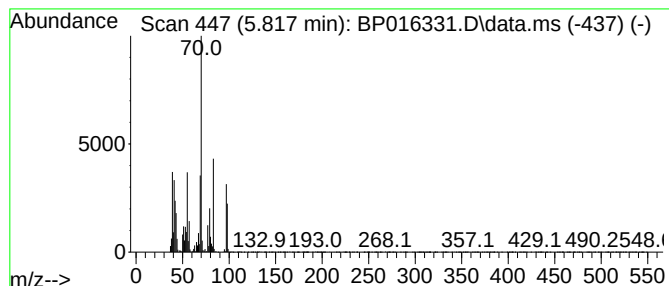
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Cyclohexen-1-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.817	15.85 ng/ul	482207	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	96
2	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	94
3	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	91
4	2-Methylene cyclopentanol			98	C6H10O	020461-31-8	83
5	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	47



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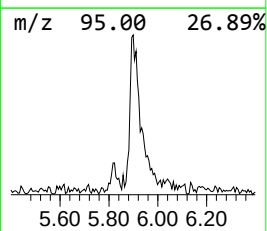
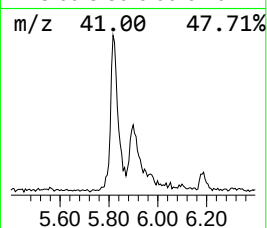
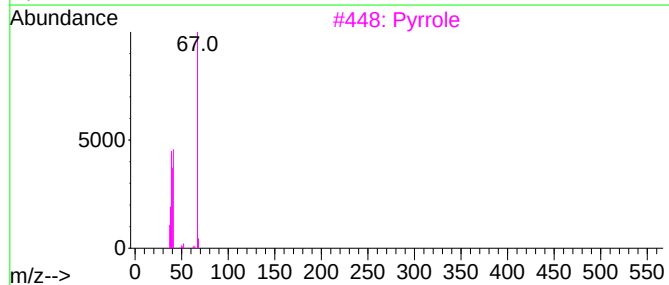
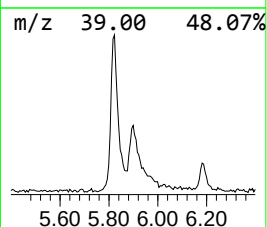
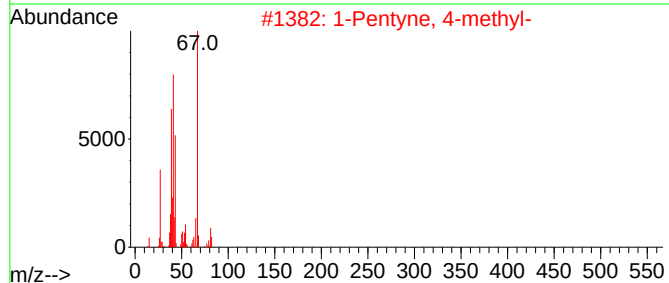
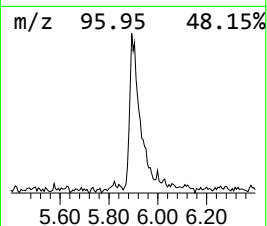
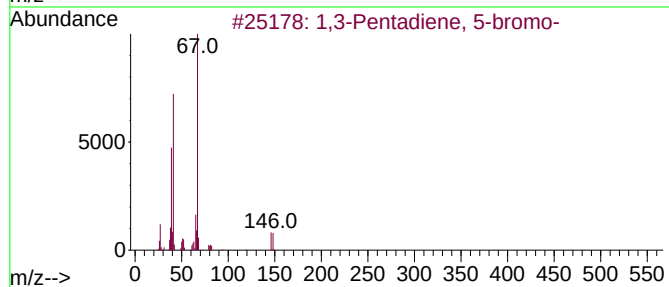
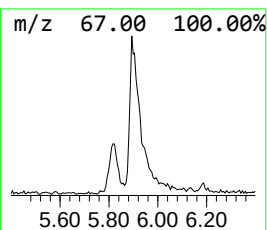
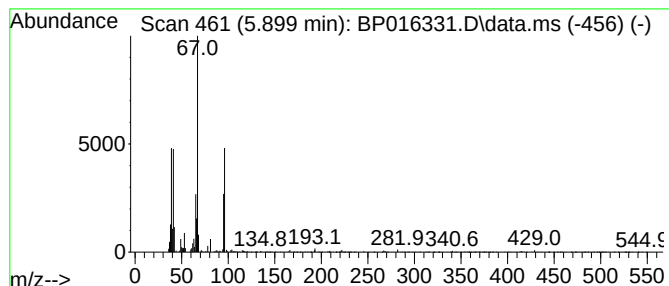
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.899	5.24 ng/ul	159418	1,4-Dichlorobenzene-d4	7.952

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3-Pentadiene, 5-bromo-	146	C5H7Br	001001-93-0	47
2		1-Pentyne, 4-methyl-	82	C6H10	007154-75-8	38
3		Pyrrole	67	C4H5N	000109-97-7	38
4		1-Cyclopentene-1-carboxylic acid	112	C6H8O2	001560-11-8	38
5		7-Bromo-4,5-diazaspiro[2.4]hept-	174	C5H7BrN2	1000261-52-7	36



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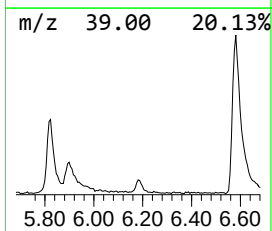
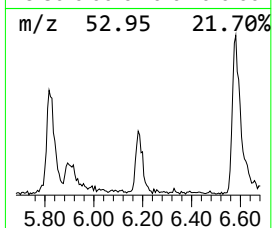
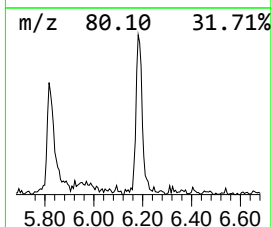
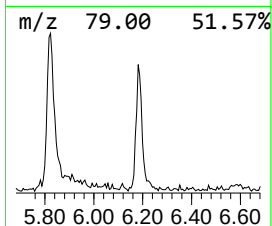
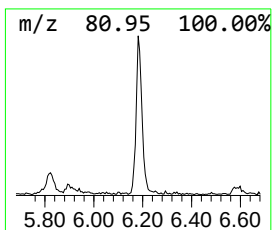
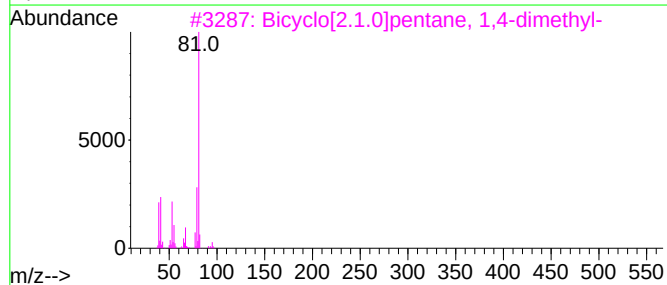
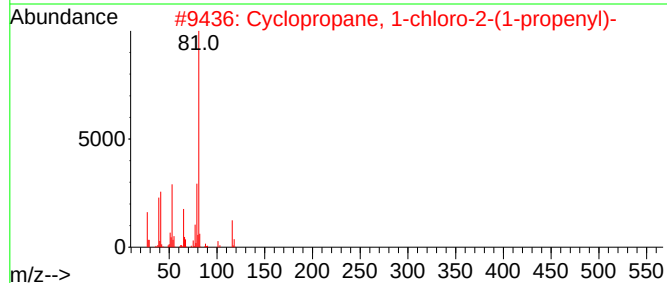
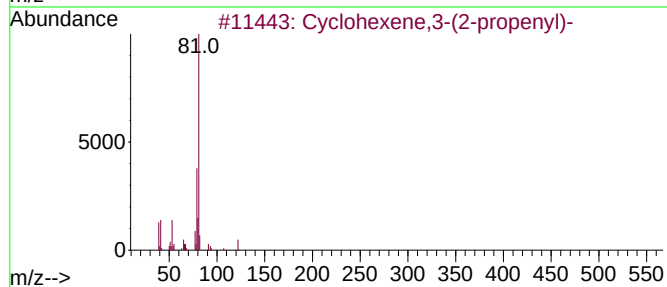
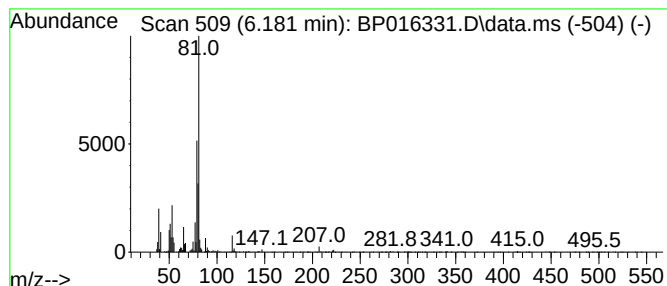
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexene,3-(2-propenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.181	2.51 ng/ul	76425	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene,3-(2-propenyl)-	122	C9H14	015232-95-8	72
2			Cyclopropane, 1-chloro-2-(1-prop...	116	C6H9Cl	074752-94-6	64
3			Bicyclo[2.1.0]pentane, 1,4-dimet...	96	C7H12	017065-18-8	53
4			Bi-2-cyclohexen-1-yl	162	C12H18	001541-20-4	53
5			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
Operator : MA/JU
Sample : 03472-11
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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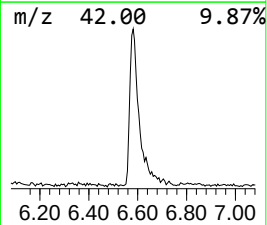
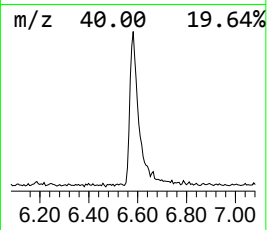
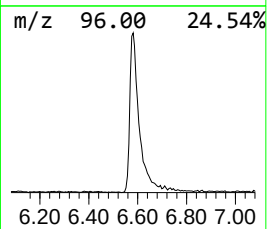
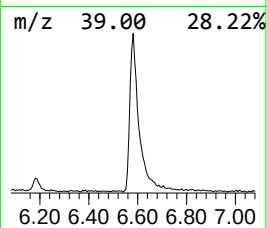
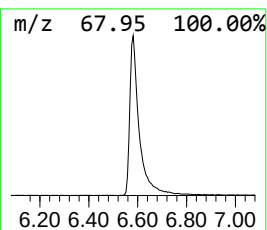
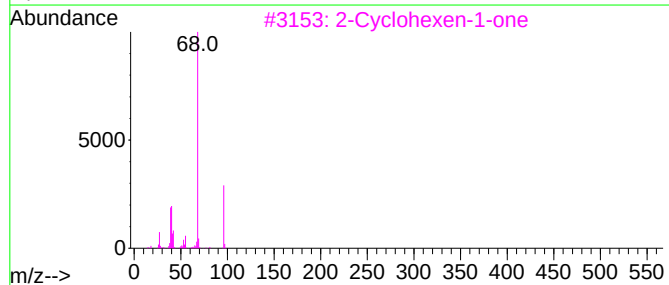
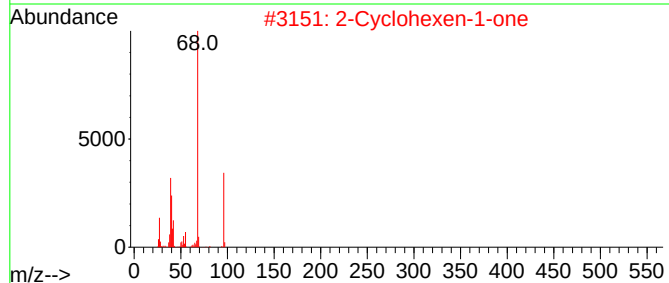
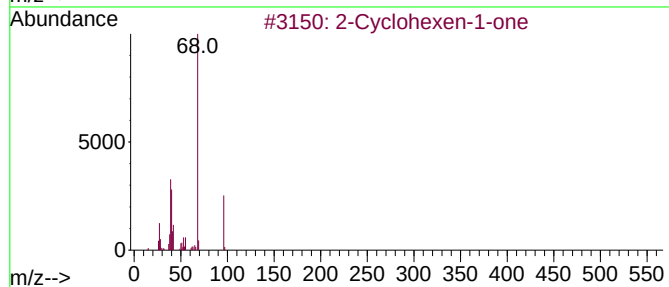
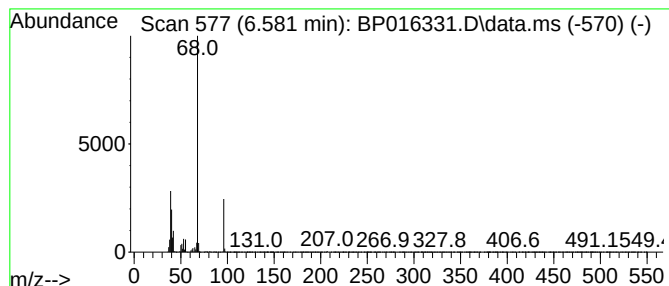
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Cyclohexen-1-one Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.581	23.23 ng/ul	706949	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	94
2			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	91
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	90
4			3-Butyn-2-amine, 2-methyl-	83	C5H9N	002978-58-7	59
5			7-Oxabicyclo[2.2.1]hept-5-en-2-one	110	C6H6O2	095530-78-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
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Instrument :
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ClientSampleId :
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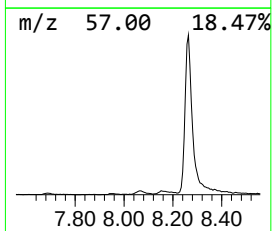
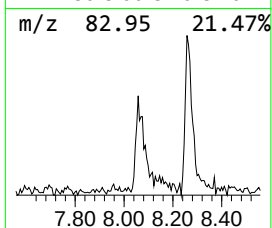
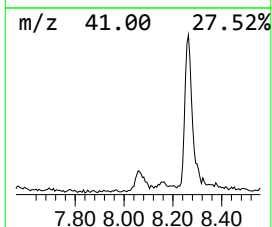
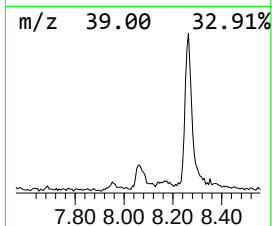
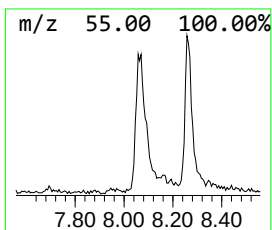
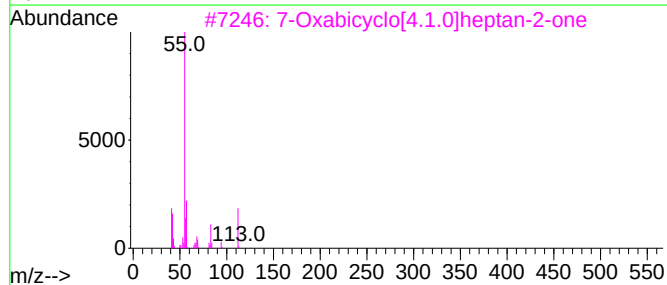
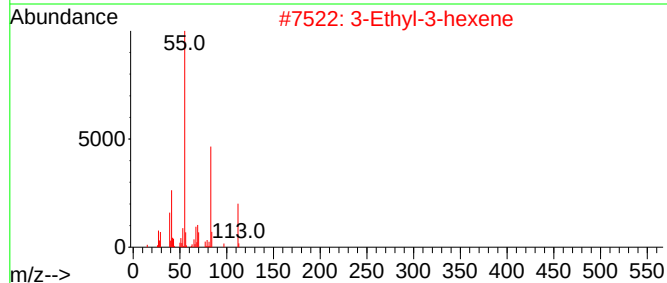
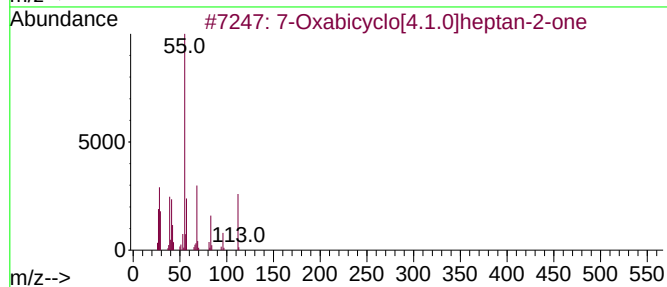
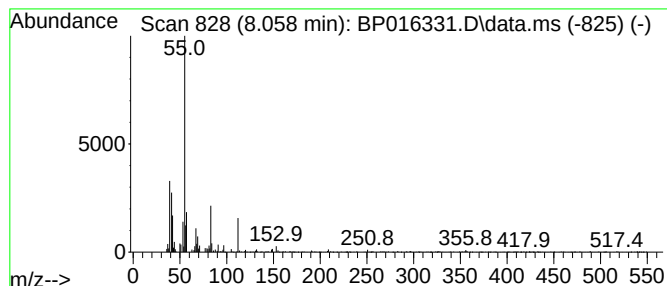
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 7-Oxabicyclo[4.1.0]heptan-2... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.058	2.94 ng/ul	89452	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[4.1.0]heptan-2-one	112	C6H8O2	006705-49-3	59
2			3-Ethyl-3-hexene	112	C8H16	016789-51-8	59
3			7-Oxabicyclo[4.1.0]heptan-2-one	112	C6H8O2	006705-49-3	49
4			2-Butenoic acid, 2-methyl-, meth...	114	C6H10O2	005953-76-4	47
5			1-Pentene, 3-ethyl-3-methyl-	112	C8H16	006196-60-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
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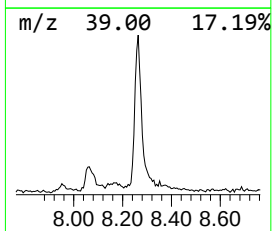
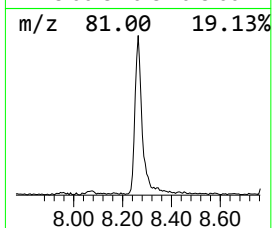
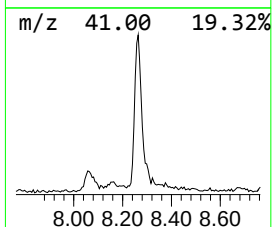
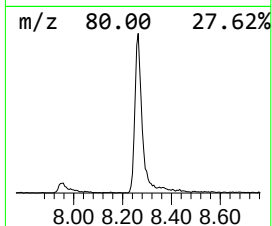
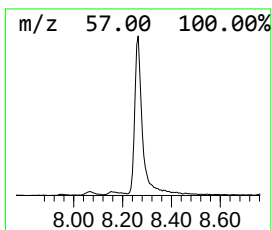
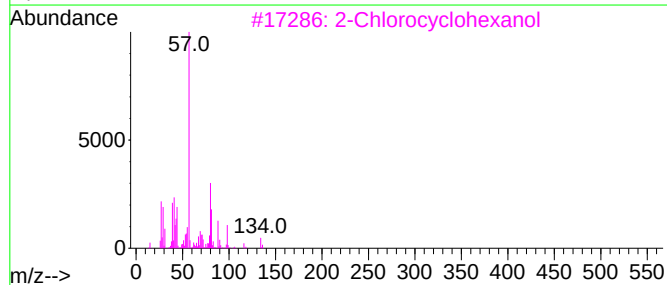
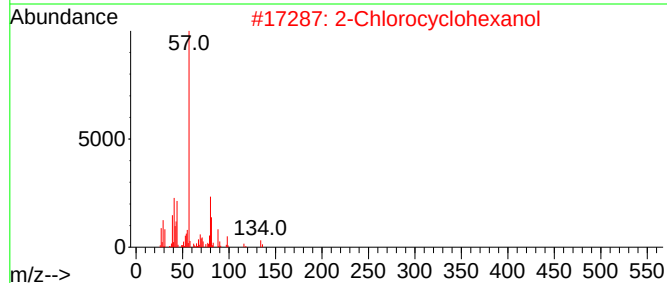
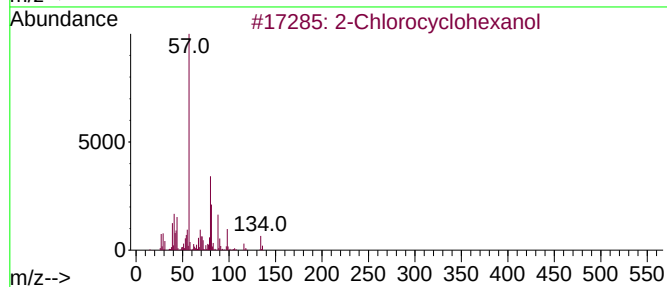
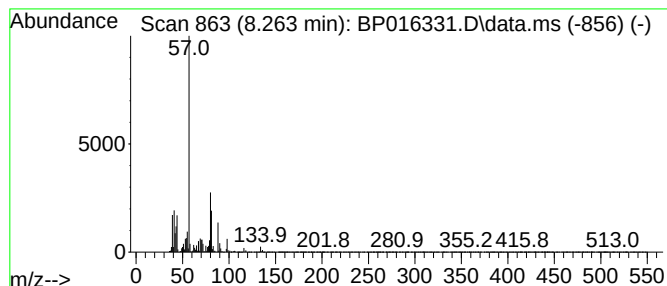
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-Chlorocyclohexanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.263	21.64 ng/ul	658514	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	91
2	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	90
3	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	64
4	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	53
5	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
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Sample : 03472-11
Misc :
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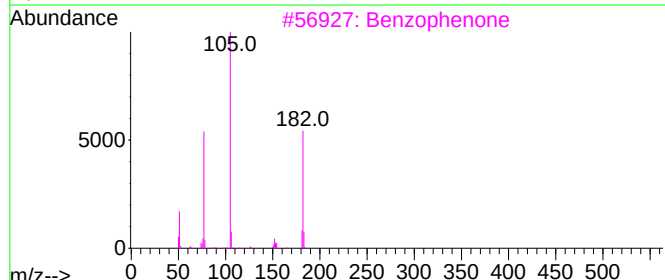
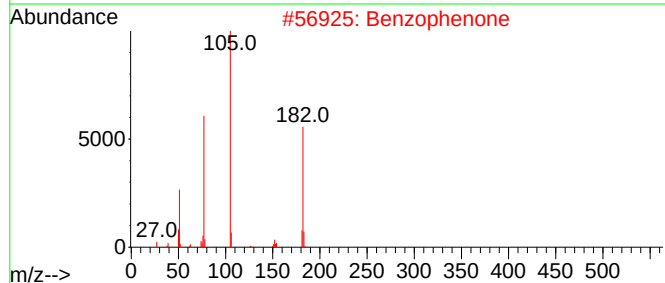
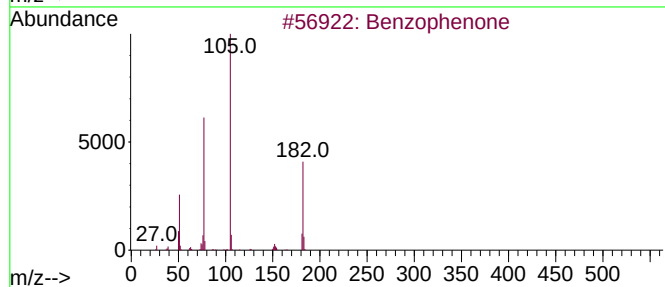
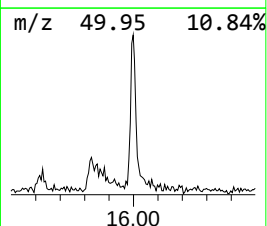
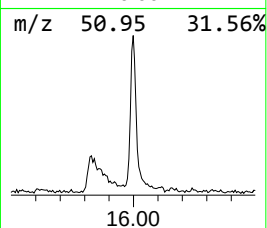
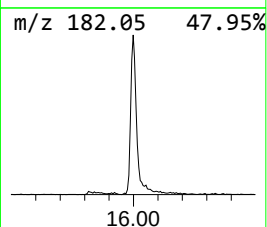
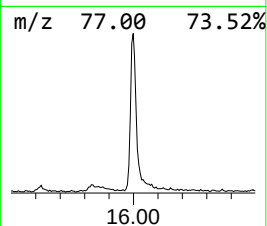
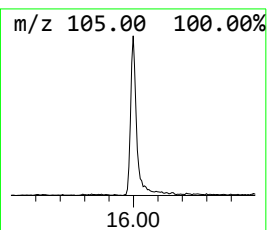
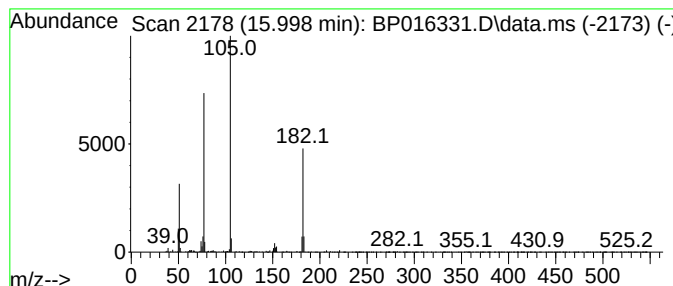
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.998	2.25 ng/ul	213649	Phenanthrene-d10	17.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
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Sample : 03472-11
Misc :
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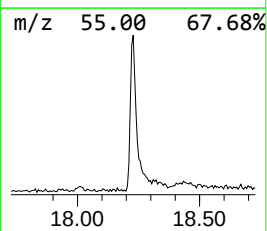
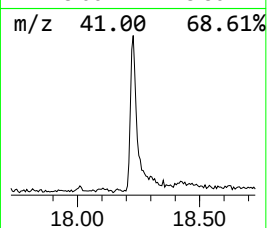
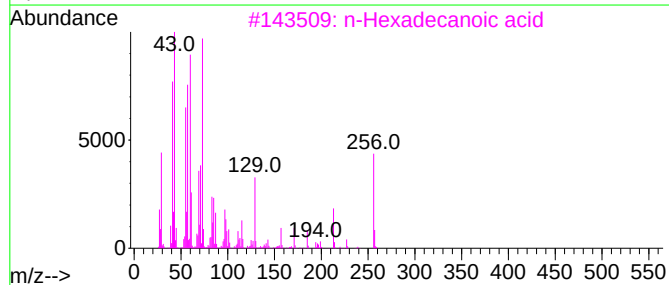
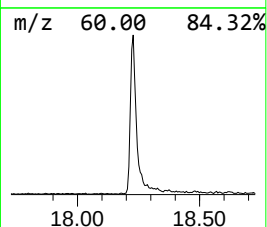
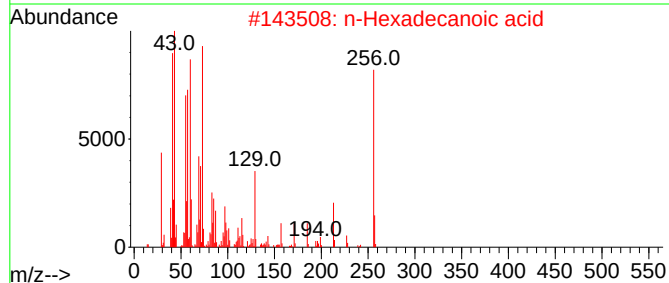
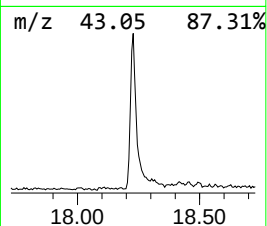
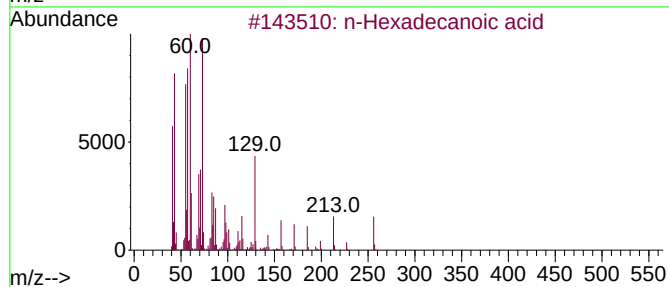
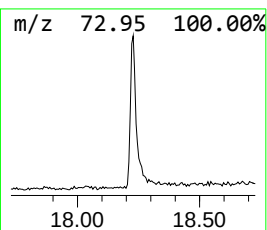
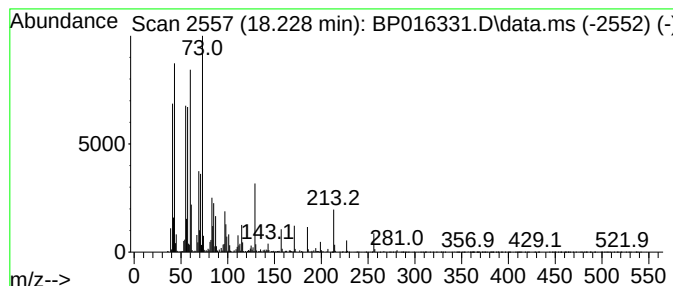
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	5.30 ng/ul	502517	Phenanthrene-d10	17.369

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
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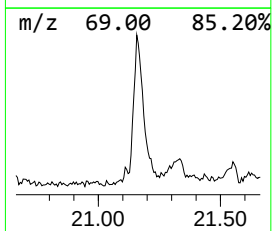
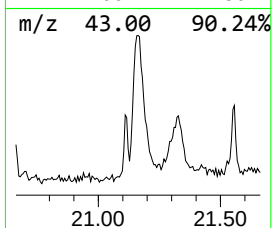
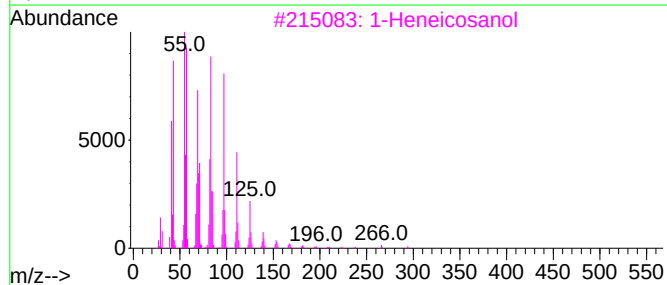
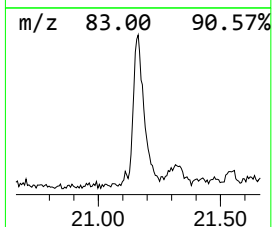
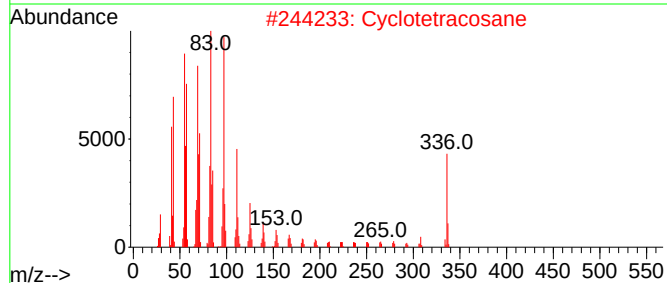
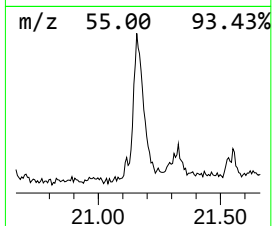
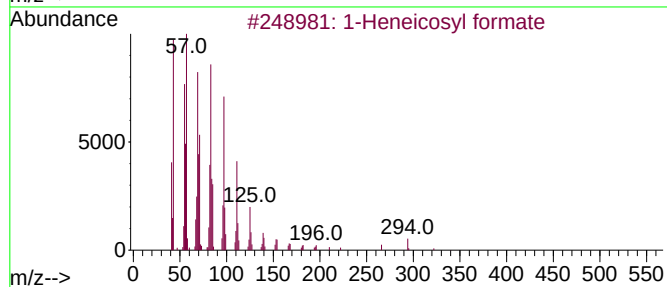
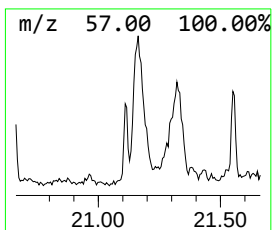
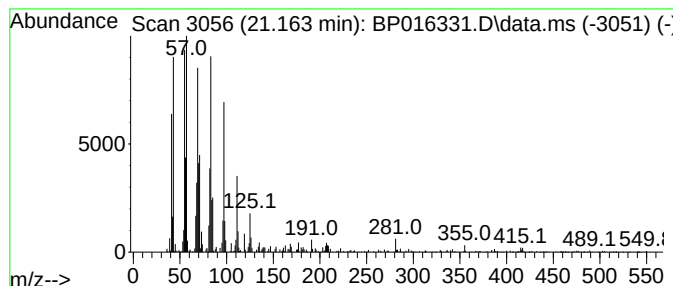
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heneicosyl formate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.163	3.74 ng/ul	453224	Chrysene-d12	21.475

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2		Cyclotetracosane	336	C24H48	000297-03-0	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	94
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5		Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071923\
Data File : BP016331.D
Acq On : 19 Jul 2023 19:20
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
Formamide, N,N-...	4.217	4.0	ng/ul	120709	1	7.952	608582	20.0
2-Cyclohexen-1-ol	5.817	15.9	ng/ul	482207	1	7.952	608582	20.0
unknown-01	5.899	5.2	ng/ul	159418	1	7.952	608582	20.0
Cyclohexene,3-(...	6.181	2.5	ng/ul	76425	1	7.952	608582	20.0
2-Cyclohexen-1-one	6.581	23.2	ng/ul	706949	1	7.952	608582	20.0
7-Oxabicyclo[4....	8.058	2.9	ng/ul	89452	1	7.952	608582	20.0
2-Chlorocyclohe...	8.263	21.6	ng/ul	658514	1	7.952	608582	20.0
Benzophenone	15.998	2.3	ng/ul	213649	4	17.369	1895620	20.0
n-Hexadecanoic ...	18.228	5.3	ng/ul	502517	4	17.369	1895620	20.0
1-Heneicosyl fo...	21.163	3.7	ng/ul	453224	5	21.475	2421520	20.0