

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
 Data File : BP015106.D
 Acq On : 16 May 2023 12:49
 Operator : CG/JU
 Sample : 02768-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONN3

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-BP051123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	5	8	14	rVB	89003	122224	0.66%	0.076%
2	3.628	36	41	53	rVB	335379	516656	2.79%	0.323%
3	3.746	55	61	70	rVB	827760	1207523	6.51%	0.754%
4	3.881	82	84	95	rBV	197120	341950	1.84%	0.214%
5	3.975	95	100	112	rVB2	302370	492371	2.66%	0.308%
6	4.181	129	135	139	rBV	225225	339704	1.83%	0.212%
7	4.752	225	232	238	rVV	8528809	13379432	72.16%	8.357%
8	4.799	238	240	245	rVV	99688	132571	0.72%	0.083%
9	5.005	265	275	278	rBV	496244	735313	3.97%	0.459%
10	5.052	278	283	293	rVV	11154093	18540977	100.00%	11.581%
11	5.134	293	297	311	rVB2	141476	224980	1.21%	0.141%
12	5.263	313	319	332	rBV	672611	1026191	5.53%	0.641%
13	5.440	342	349	354	rBV	149850	231383	1.25%	0.145%
14	6.357	499	505	514	rVB	1254718	1929358	10.41%	1.205%
15	7.451	685	691	698	rVV	1409378	2157243	11.64%	1.347%
16	7.534	698	705	717	rVB	1663864	2573657	13.88%	1.608%
17	7.657	719	726	737	rBV	900824	1447119	7.80%	0.904%
18	7.840	748	757	767	rVB	976668	1540495	8.31%	0.962%
19	8.134	799	807	812	rBV	1248009	1985831	10.71%	1.240%
20	8.193	812	817	828	rVB	828350	1315168	7.09%	0.821%
21	8.393	845	851	860	rBV2	103540	177485	0.96%	0.111%
22	8.504	863	870	880	rVB2	358119	589805	3.18%	0.368%
23	8.634	886	892	899	rVB	304547	479056	2.58%	0.299%
24	8.740	901	910	922	rBV	179030	338425	1.83%	0.211%
25	8.846	922	928	939	rVV	62482	125140	0.67%	0.078%
26	9.310	991	1007	1018	rBV	1653392	3037844	16.38%	1.897%
27	9.416	1018	1025	1029	rVV	863960	1696805	9.15%	1.060%
28	9.504	1029	1040	1055	rVB	1444422	4512194	24.34%	2.818%
29	9.704	1069	1074	1080	rVB	125505	190503	1.03%	0.119%
30	10.040	1122	1131	1146	rBV	1118115	1898116	10.24%	1.186%
31	10.187	1151	1156	1164	rBV	69411	126831	0.68%	0.079%
32	10.575	1215	1222	1236	rBV	2478601	4270365	23.03%	2.667%
33	10.722	1240	1247	1258	rVV2	218634	540815	2.92%	0.338%
34	10.822	1258	1264	1278	rVV	260141	474747	2.56%	0.297%
35	10.963	1280	1288	1297	rVB	1738206	2919217	15.74%	1.823%
36	11.116	1306	1314	1322	rVV	368701	611175	3.30%	0.382%

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37	11.892	1441	1446	1456	rVV2	130263	253129	1.37%	0.158%
38	11.992	1457	1463	1469	rVV	313297	560873	3.03%	0.350%
39	12.204	1489	1499	1508	rBV3	117994	384803	2.08%	0.240%
40	12.416	1528	1535	1545	rVV	116220	203389	1.10%	0.127%
41	12.639	1567	1573	1588	rBV	604188	991092	5.35%	0.619%
42	12.904	1609	1618	1629	rBV3	61309	125126	0.67%	0.078%
43	13.192	1655	1667	1678	rBV	997918	1710139	9.22%	1.068%
44	13.757	1754	1763	1770	rBV	185987	291601	1.57%	0.182%
45	14.175	1823	1834	1839	rBV	4300461	6111921	32.96%	3.818%
46	14.216	1839	1841	1846	rVV	102717	139180	0.75%	0.087%
47	14.269	1846	1850	1856	rVV	303254	459521	2.48%	0.287%
48	14.328	1856	1860	1870	rVB	121374	187911	1.01%	0.117%
49	14.469	1878	1884	1894	rVV	1167417	1655410	8.93%	1.034%
50	14.669	1913	1918	1928	rVB2	168798	258120	1.39%	0.161%
51	14.775	1928	1936	1944	rBV	2802461	4125842	22.25%	2.577%
52	14.957	1962	1967	1979	rBV	264418	478884	2.58%	0.299%
53	15.475	2051	2055	2066	rBV	48537	119493	0.64%	0.075%
54	15.586	2070	2074	2086	rVV4	82681	177825	0.96%	0.111%
55	15.763	2098	2104	2117	rVV	5119356	7503940	40.47%	4.687%
56	15.880	2118	2124	2138	rVV	1826541	2526637	13.63%	1.578%
57	16.127	2160	2166	2177	rVB	185328	270233	1.46%	0.169%
58	16.457	2216	2222	2231	rBV3	105575	199424	1.08%	0.125%
59	17.180	2337	2345	2352	rBV	606387	863373	4.66%	0.539%
60	17.398	2377	2382	2390	rBV2	86772	139884	0.75%	0.087%
61	17.527	2397	2404	2412	rBV	3572128	4964151	26.77%	3.101%
62	17.622	2412	2420	2431	rVV2	3314450	5758418	31.06%	3.597%
63	17.710	2431	2435	2439	rVV	174349	271002	1.46%	0.169%
64	17.763	2439	2444	2459	rVB	1287062	2182702	11.77%	1.363%
65	18.327	2535	2540	2545	rVV2	280991	385009	2.08%	0.240%
66	18.380	2545	2549	2554	rVV	264380	375844	2.03%	0.235%
67	18.457	2559	2562	2570	rVB	71050	118639	0.64%	0.074%
68	18.563	2576	2580	2586	rVB	83904	113449	0.61%	0.071%
69	18.869	2627	2632	2641	rBV2	196943	298247	1.61%	0.186%
70	19.127	2671	2676	2680	rBV	968732	1423843	7.68%	0.889%
71	19.169	2680	2683	2695	rVB3	342947	618329	3.33%	0.386%
72	19.486	2732	2737	2746	rBV2	268464	514348	2.77%	0.321%
73	19.604	2753	2757	2765	rBV2	131411	188763	1.02%	0.118%
74	19.845	2792	2798	2812	rVB	5452879	7033521	37.94%	4.393%
75	20.263	2864	2869	2875	rBV2	324692	438527	2.37%	0.274%
76	20.386	2887	2890	2894	rVB2	101852	118854	0.64%	0.074%

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 Title : SVOA CALIBRATION

77	20.604	2924	2927	2929	rBV2	81849	113514	0.61%	0.071%
78	20.639	2930	2933	2937	rVB	244239	275014	1.48%	0.172%
79	20.833	2962	2966	2969	rBV3	297904	447221	2.41%	0.279%
80	20.880	2969	2974	2975	rVV2	337200	501866	2.71%	0.313%
81	20.898	2975	2977	2982	rVV	831369	944295	5.09%	0.590%
82	20.945	2982	2985	2988	rVV	314492	318398	1.72%	0.199%
83	21.021	2994	2998	3003	rBV	1104466	1406508	7.59%	0.879%
84	21.086	3003	3009	3013	rBV	615047	1125254	6.07%	0.703%
85	21.133	3013	3017	3019	rBV	590045	807778	4.36%	0.505%
86	21.198	3025	3028	3032	rVB	511793	621837	3.35%	0.388%
87	21.245	3033	3036	3039	rBV4	101889	143104	0.77%	0.089%
88	21.286	3039	3043	3046	rBV2	214821	347446	1.87%	0.217%
89	21.604	3092	3097	3107	rVB	3734173	5053702	27.26%	3.157%
90	21.810	3129	3132	3134	rBV2	183357	255879	1.38%	0.160%
91	21.886	3141	3145	3151	rVB	369014	573430	3.09%	0.358%
92	21.951	3152	3156	3158	rBV3	429511	672100	3.62%	0.420%
93	22.104	3177	3182	3184	rBV2	745777	1185933	6.40%	0.741%
94	22.198	3194	3198	3207	rVB2	513325	1238635	6.68%	0.774%
95	22.480	3241	3246	3259	rBV	1111853	2955395	15.94%	1.846%
96	22.757	3289	3293	3298	rBV	243138	445734	2.40%	0.278%
97	23.609	3433	3438	3448	rBV2	459204	1376185	7.42%	0.860%
98	24.004	3499	3505	3516	rVB2	1423094	3162395	17.06%	1.975%
99	24.174	3527	3534	3555	rVB	2245086	5230599	28.21%	3.267%
100	26.086	3851	3859	3872	rVB2	1444537	4129720	22.27%	2.579%

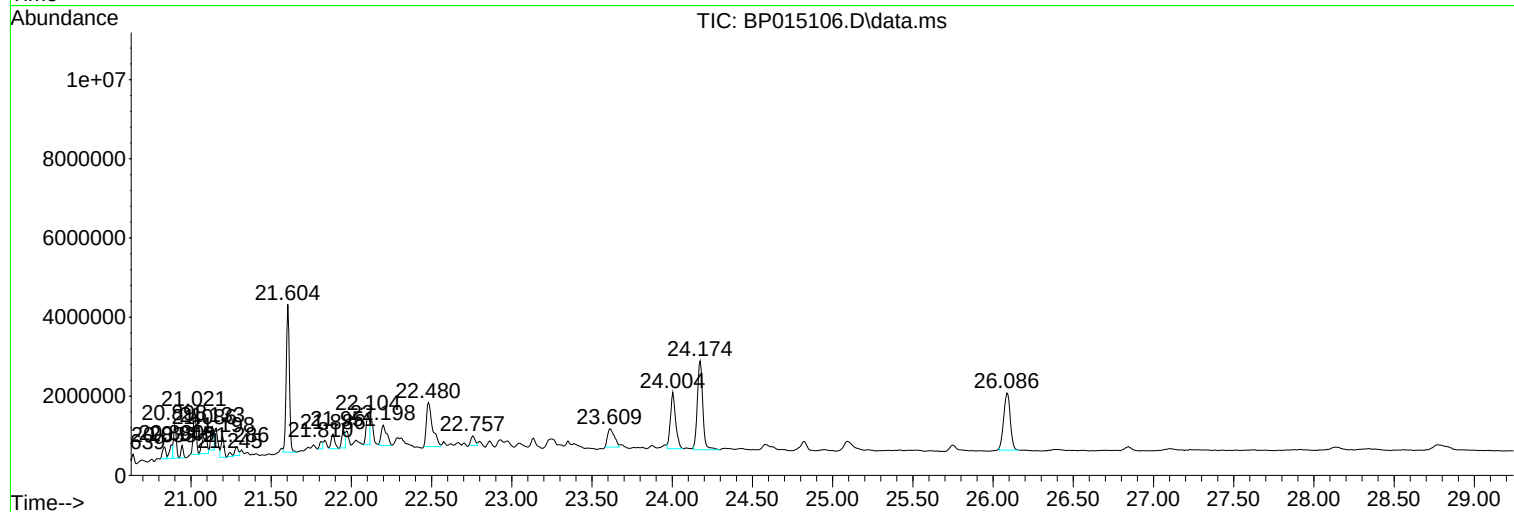
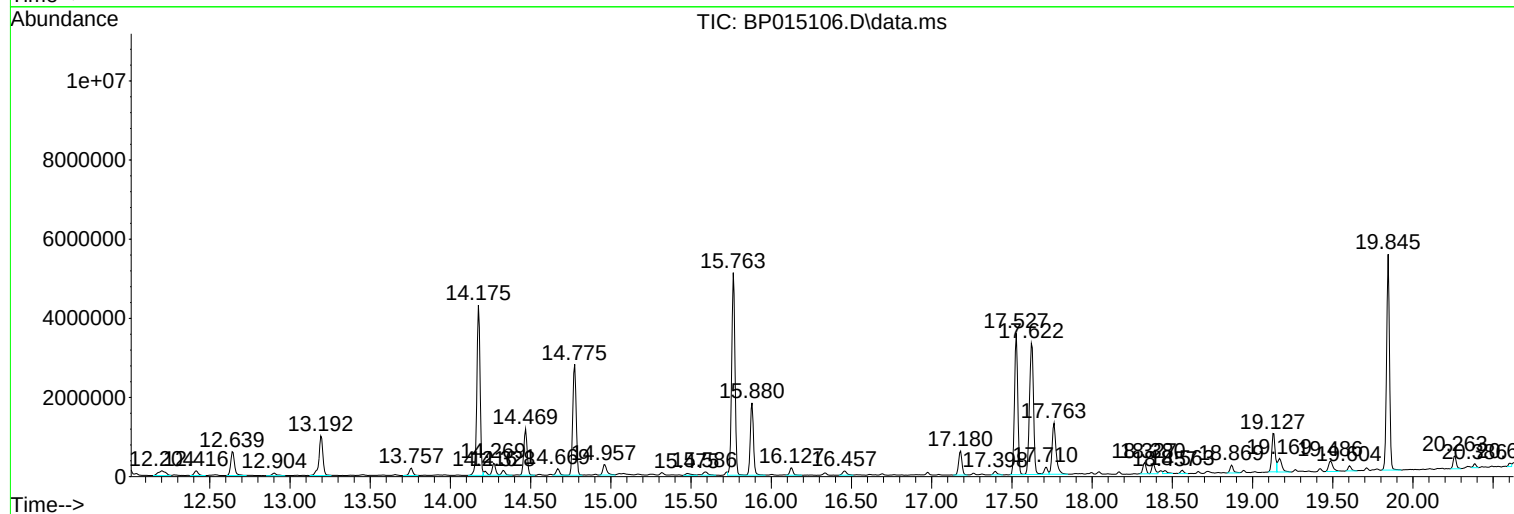
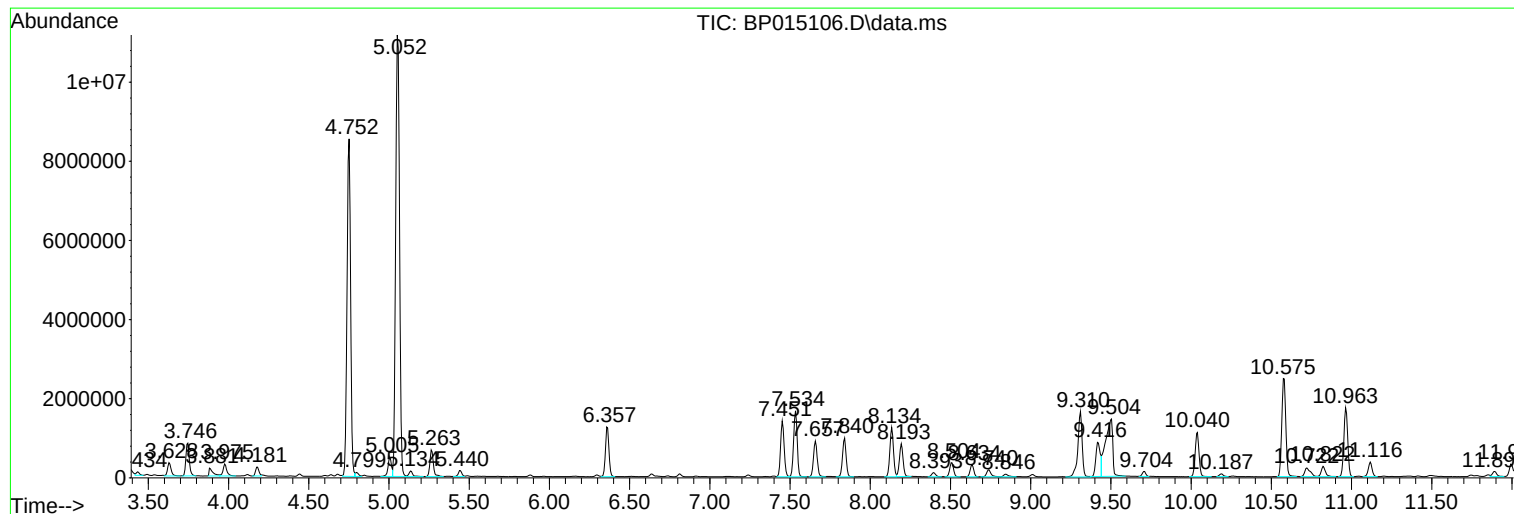
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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



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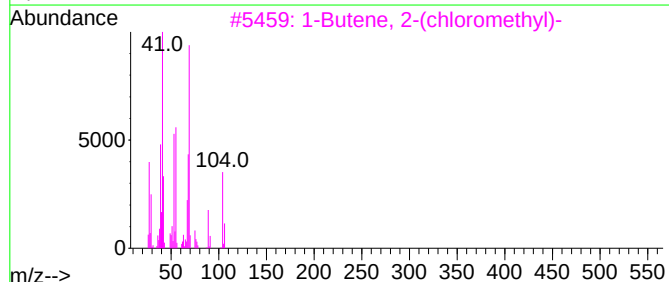
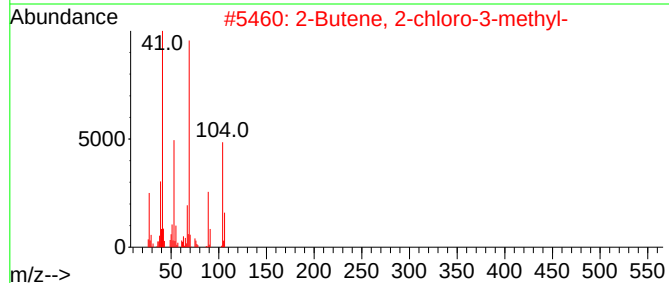
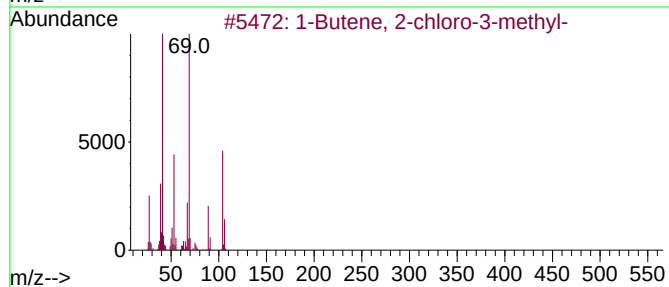
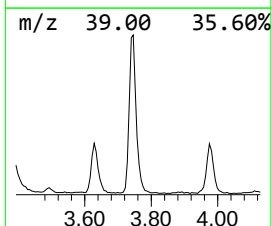
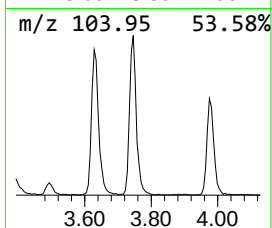
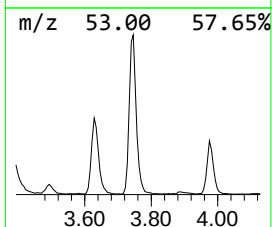
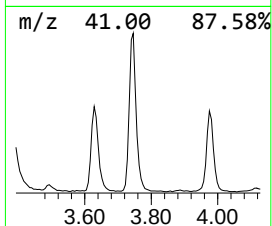
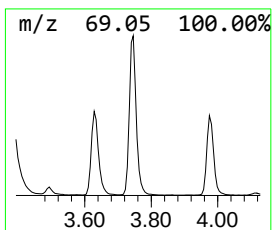
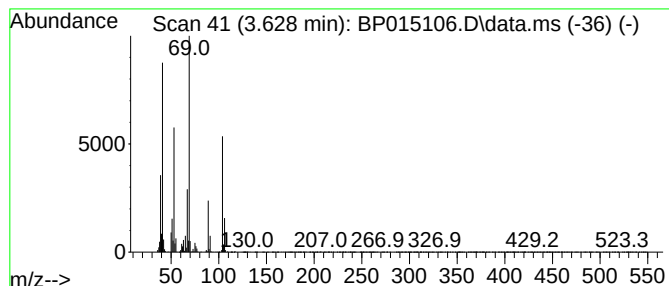
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Butene, 2-chloro-3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.628	5.20 ng/ul	516656	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butene, 2-chloro-3-methyl-		104	C5H9Cl	017773-64-7	91	
2	2-Butene, 2-chloro-3-methyl-		104	C5H9Cl	017773-65-8	90	
3	1-Butene, 2-(chloromethyl)-		104	C5H9Cl	023010-02-8	90	
4	1-Butene, 1-chloro-2-methyl-		104	C5H9Cl	023378-11-2	64	
5	1-Butene, 2-chloro-3-methyl-		104	C5H9Cl	017773-64-7	59	



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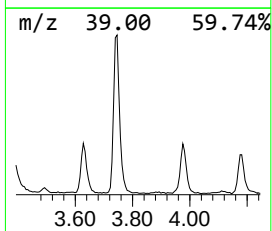
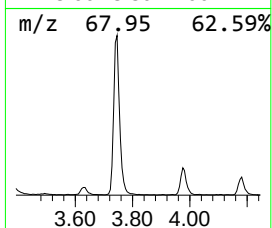
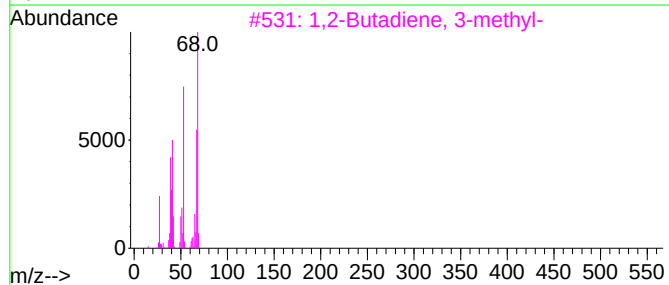
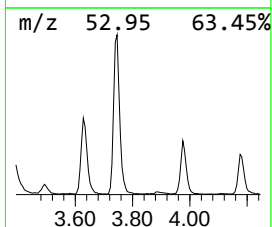
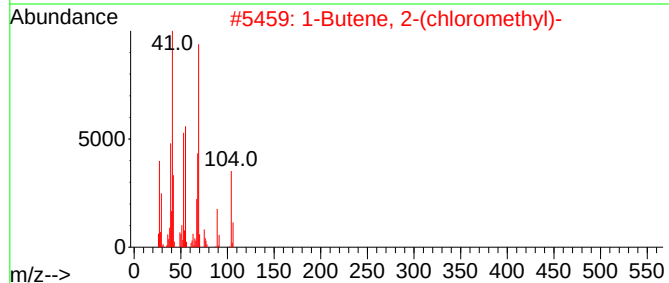
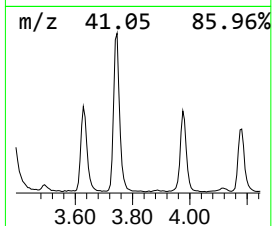
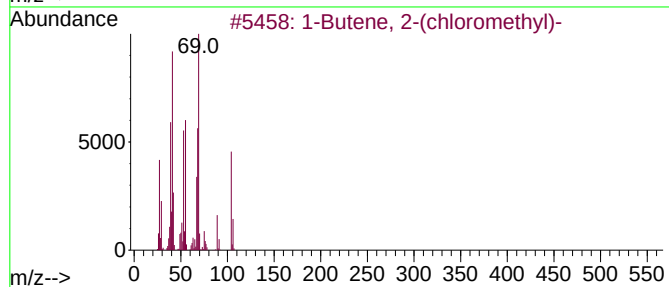
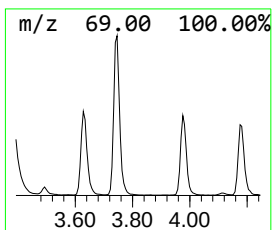
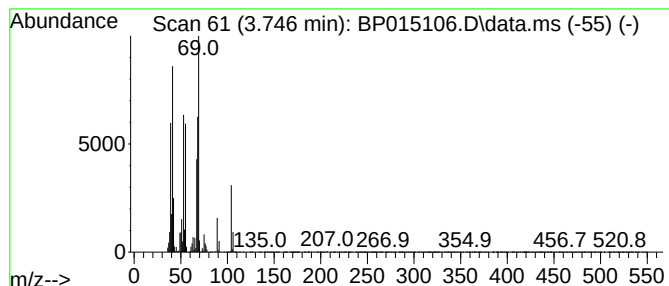
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Butene, 2-(chloromethyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.746	12.16 ng/ul	1207520	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butene, 2-(chloromethyl)-		104	C5H9Cl	023010-02-8	96	
2	1-Butene, 2-(chloromethyl)-		104	C5H9Cl	023010-02-8	91	
3	1,2-Butadiene, 3-methyl-		68	C5H8	000598-25-4	60	
4	1,2-Butadiene, 3-methyl-		68	C5H8	000598-25-4	55	
5	2-Pentyne		68	C5H8	000627-21-4	45	



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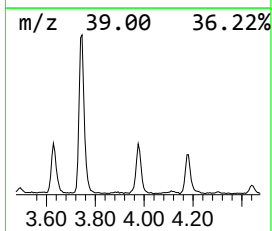
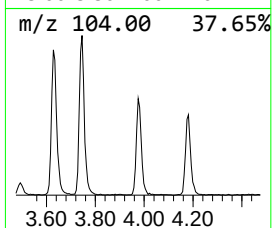
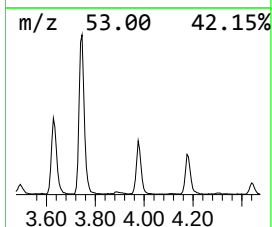
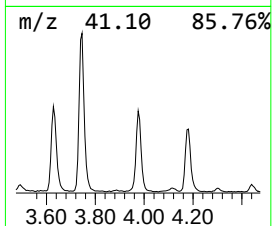
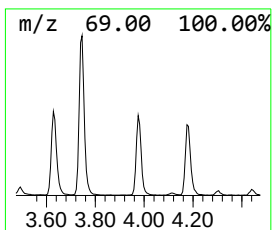
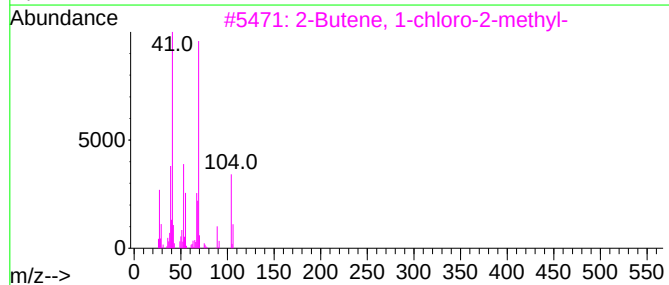
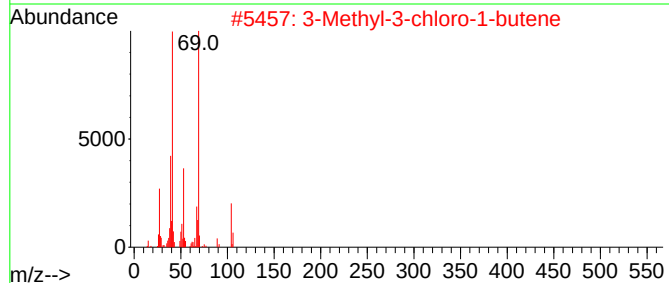
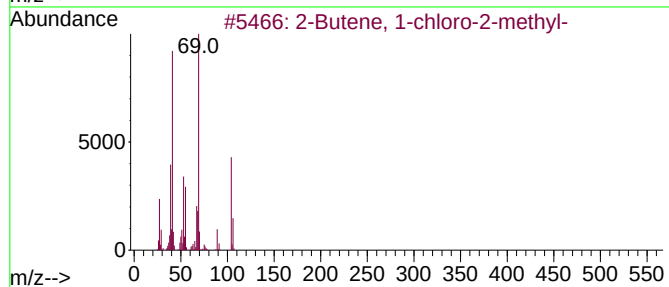
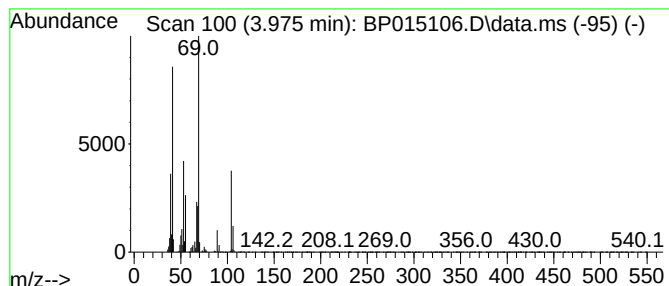
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Butene, 1-chloro-2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.975	4.96 ng/ul	492371	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 1-chloro-2-methyl-		104	C5H9Cl	013417-43-1	91	
2	3-Methyl-3-chloro-1-butene		104	C5H9Cl	002190-48-9	91	
3	2-Butene, 1-chloro-2-methyl-		104	C5H9Cl	013417-43-1	91	
4	1-Butene, 2-chloro-3-methyl-		104	C5H9Cl	017773-64-7	87	
5	1-Butene, 2-(chloromethyl)-		104	C5H9Cl	023010-02-8	78	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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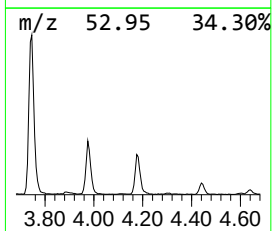
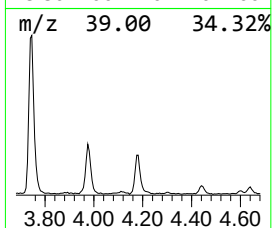
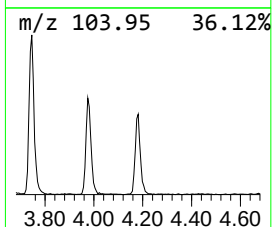
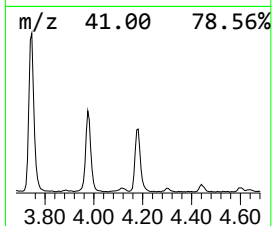
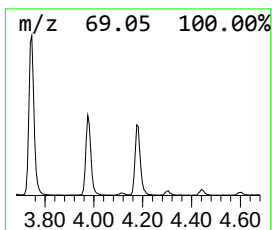
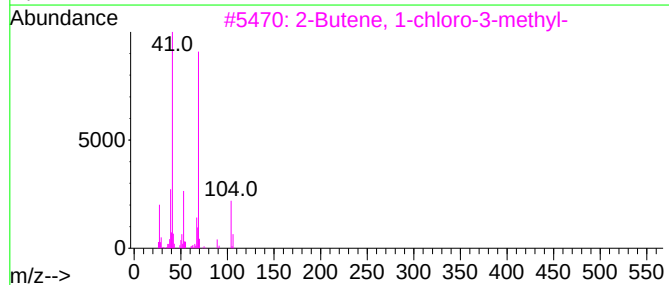
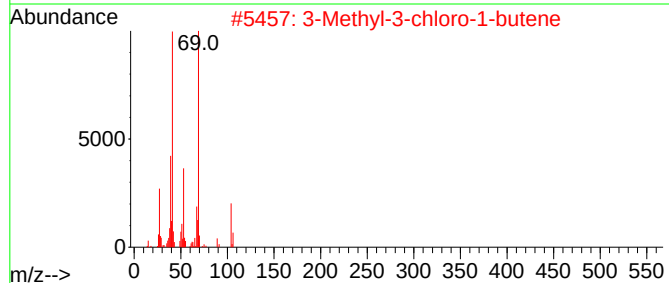
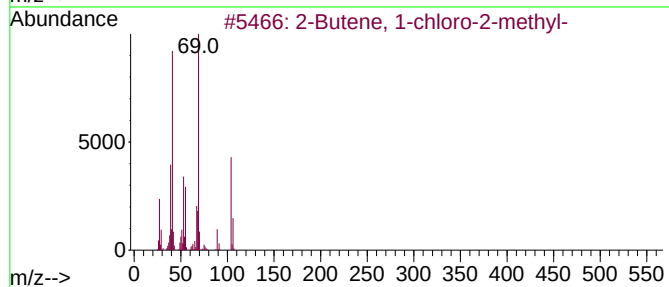
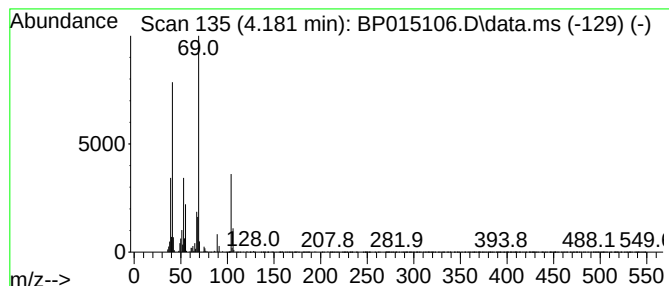
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 3-Methyl-3-chloro-1-butene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.181	3.42 ng/ul	339704	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 1-chloro-2-methyl-			104	C5H9Cl	013417-43-1	96
2	3-Methyl-3-chloro-1-butene			104	C5H9Cl	002190-48-9	91
3	2-Butene, 1-chloro-3-methyl-			104	C5H9Cl	000503-60-6	91
4	1-Butene, 2-chloro-3-methyl-			104	C5H9Cl	017773-64-7	90
5	2-Butene, 1-chloro-3-methyl-			104	C5H9Cl	000503-60-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
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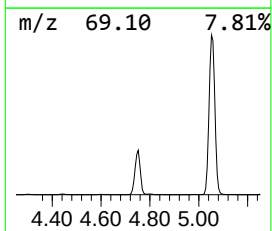
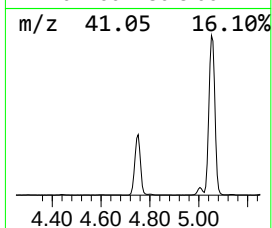
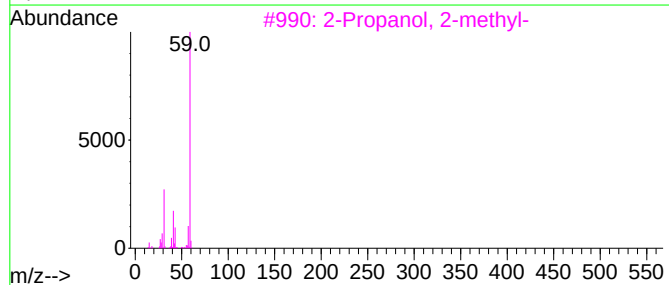
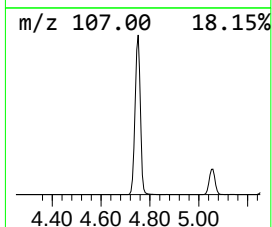
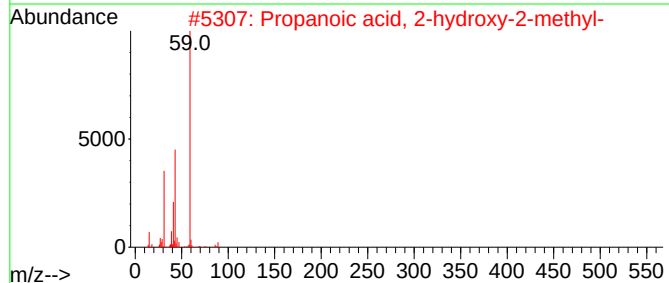
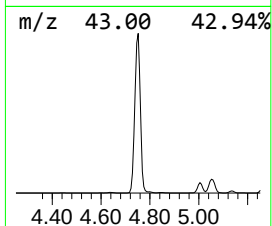
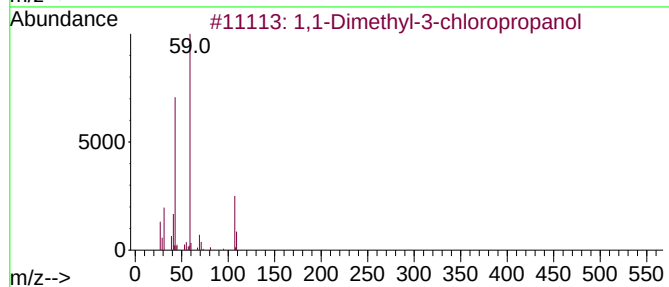
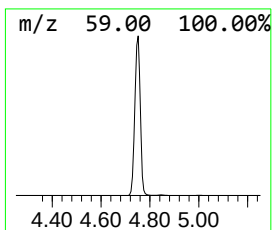
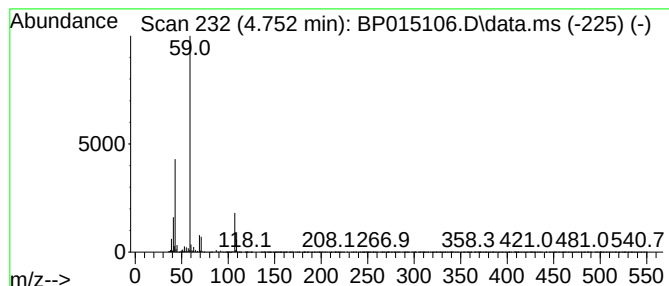
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1-Dimethyl-3-chloropropanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.752	134.75 ng/ul	13379400	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Dimethyl-3-chloropropanol	122	C5H11ClO	001985-88-2	74
2			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	59
3			2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	40
4			4-Methyl-2-pentanol, methyl ether	116	C7H16O	1000333-93-5	38
5			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	38



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Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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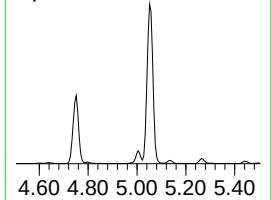
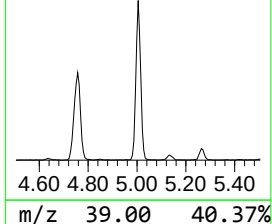
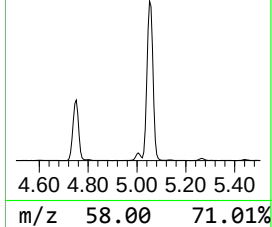
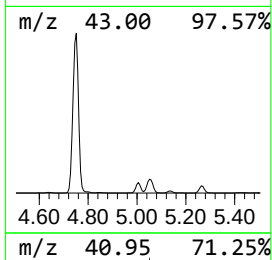
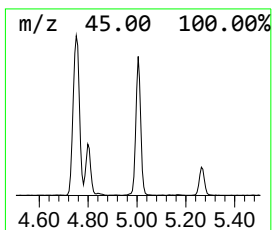
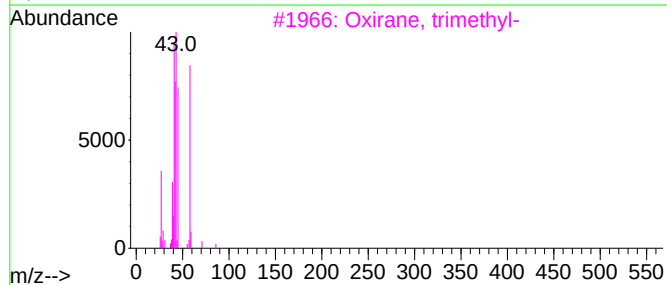
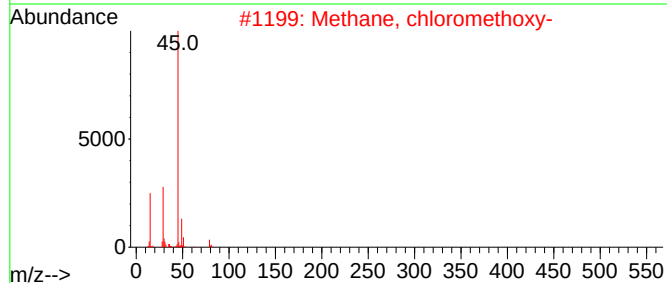
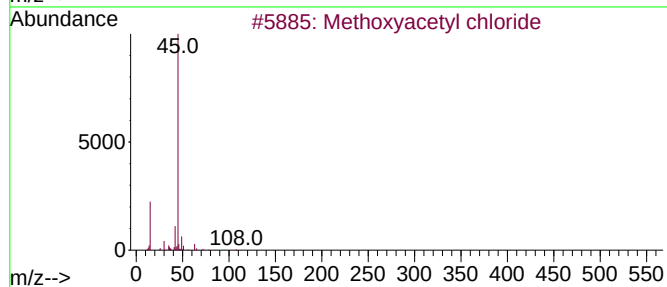
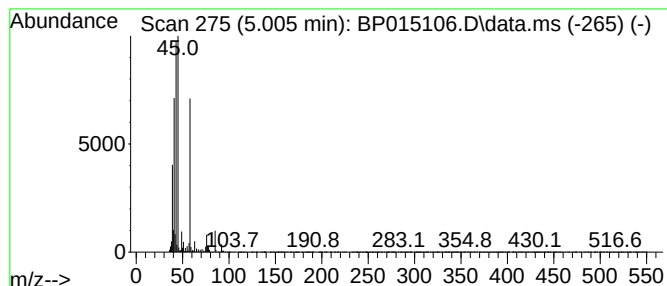
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.005	7.41 ng/ul	735313	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxyacetyl chloride	108	C3H5ClO2	038870-89-2	35
2			Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	35
3			Oxirane, trimethyl-	86	C5H10O	005076-19-7	25
4			Oxirane, trimethyl-	86	C5H10O	005076-19-7	25
5			Ethanol, 2-methoxy-, acetate	118	C5H10O3	000110-49-6	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
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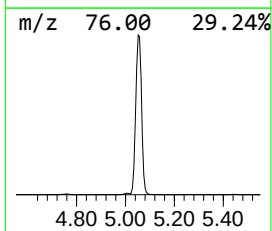
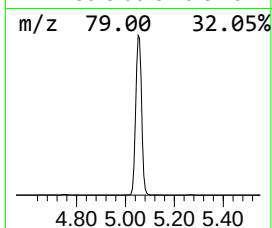
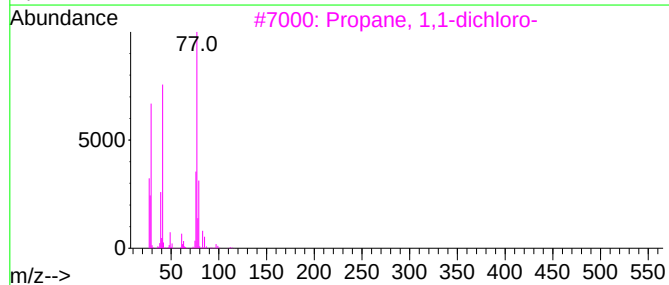
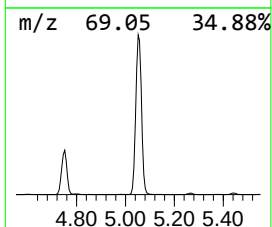
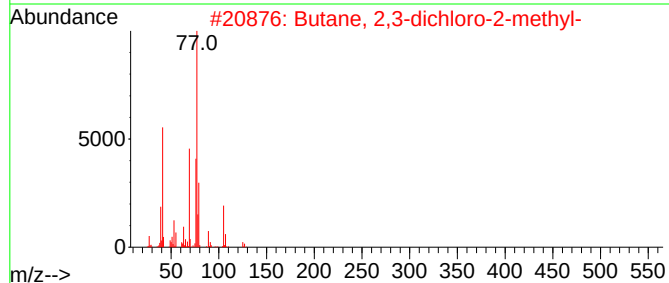
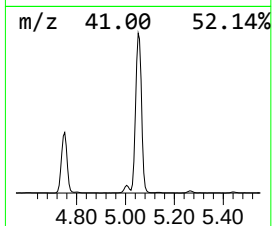
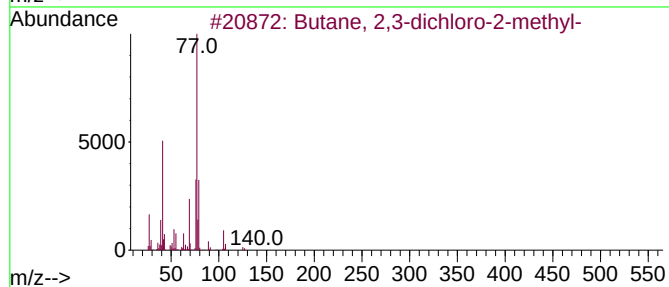
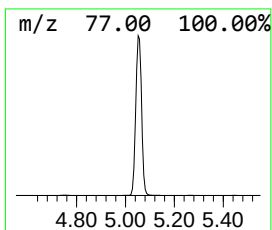
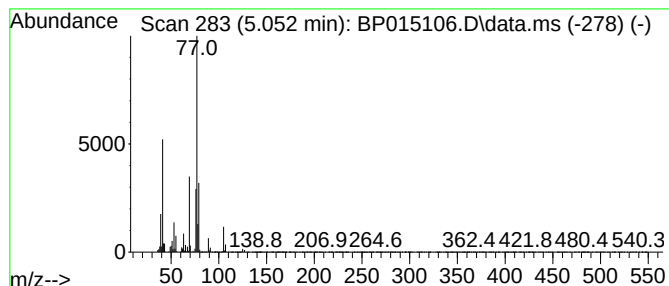
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Butane, 2,3-dichloro-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.052	186.73 ng/ul	18541000	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	90
2			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	58
3			Propane, 1,1-dichloro-	112	C3H6Cl2	000078-99-9	45
4			Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	33
5			1-Propene, 2-chloro-	76	C3H5Cl	000557-98-2	10



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Data File : BP015106.D
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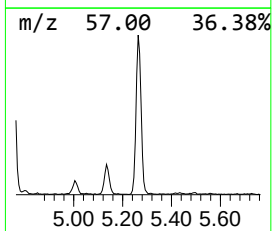
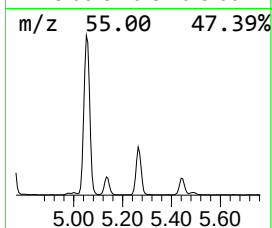
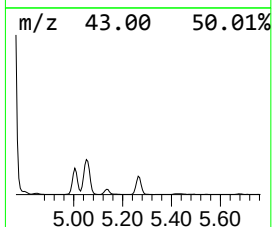
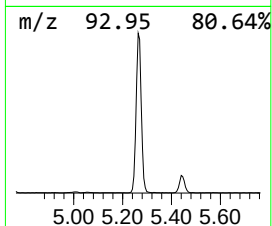
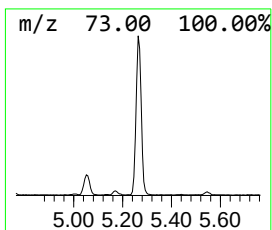
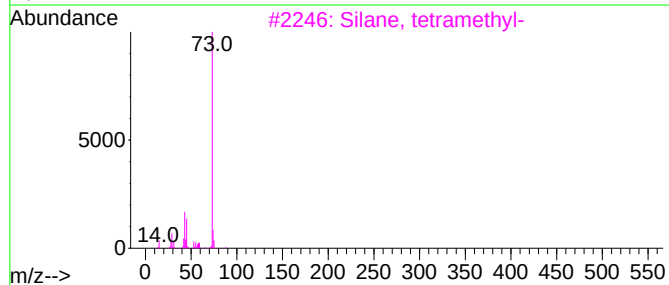
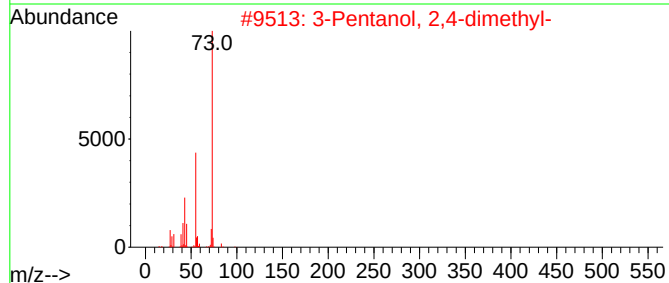
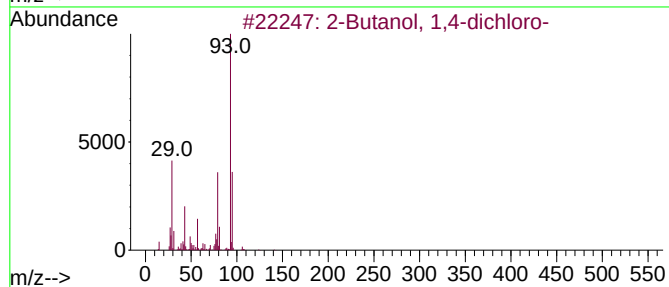
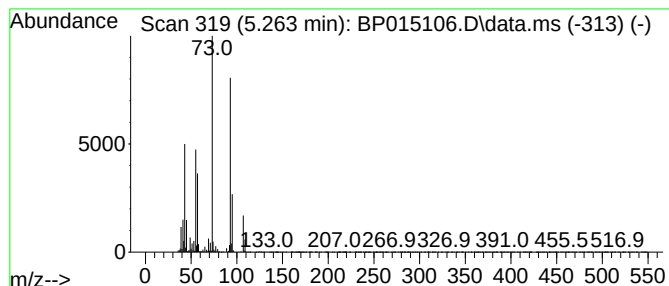
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.263	10.34 ng/ul	1026190	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 1,4-dichloro-		142	C4H8Cl2O	002419-74-1	40	
2	3-Pentanol, 2,4-dimethyl-		116	C7H16O	000600-36-2	9	
3	Silane, tetramethyl-		88	C4H12Si	000075-76-3	9	
4	Chlorodimethylethylsilane		122	C4H11ClSi	006917-76-6	9	
5	Butanoic acid, 2-hydroxy-2-methy...		132	C6H12O3	032793-34-3	9	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
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ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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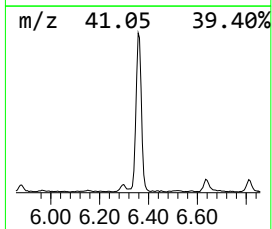
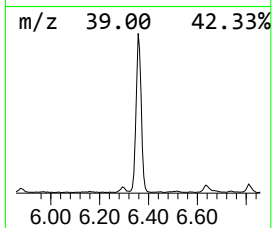
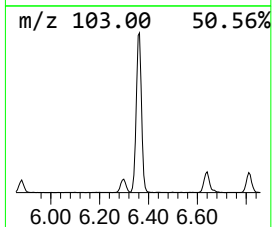
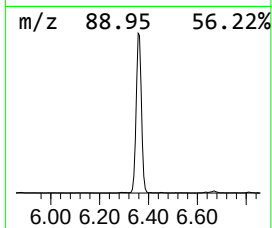
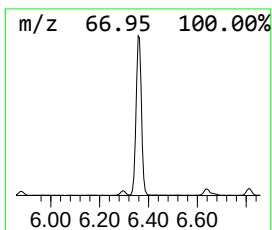
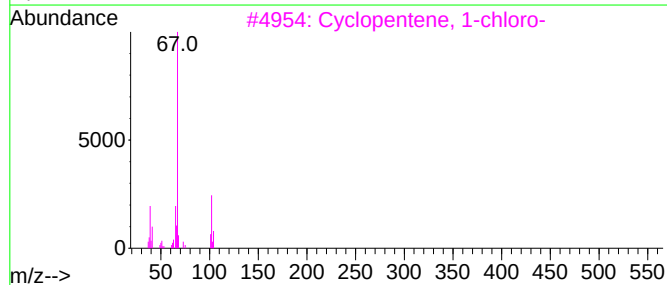
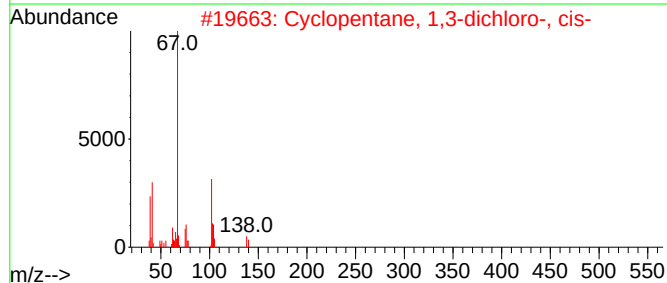
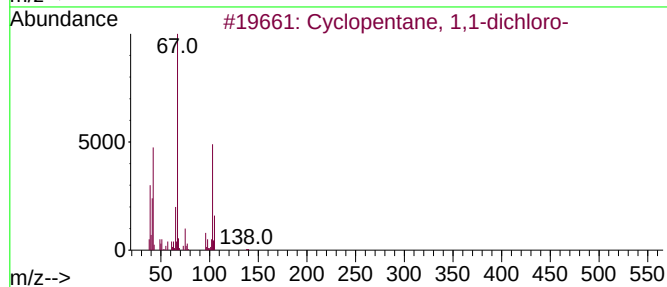
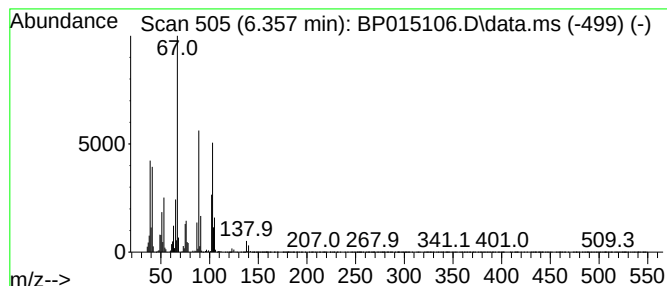
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.357	19.43 ng/ul	1929360	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1-dichloro-	138	C5H8Cl2	031038-06-9	47
2			Cyclopentane, 1,3-dichloro-, cis-	138	C5H8Cl2	026688-51-7	45
3			Cyclopentene, 1-chloro-	102	C5H7Cl	000930-29-0	43
4			1-Butyne, 3-chloro-3-methyl-	102	C5H7Cl	001111-97-3	38
5			1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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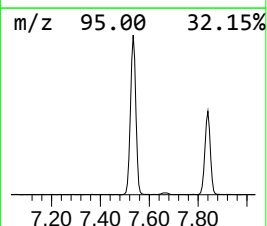
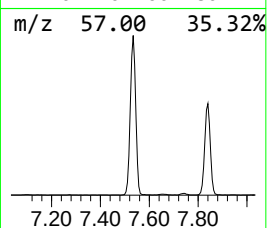
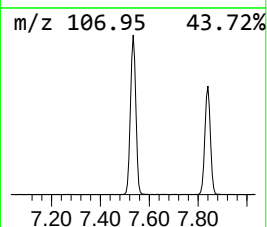
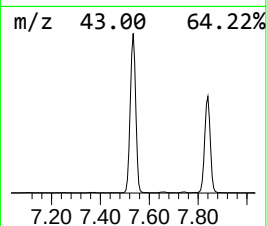
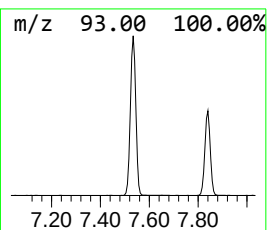
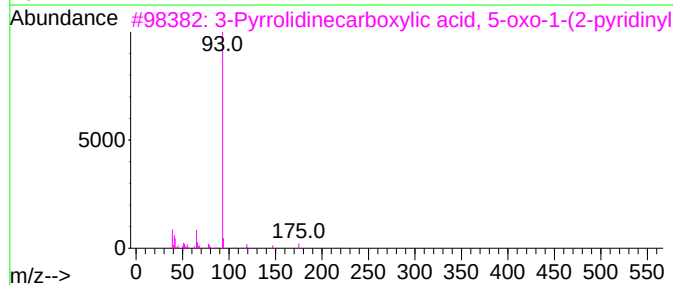
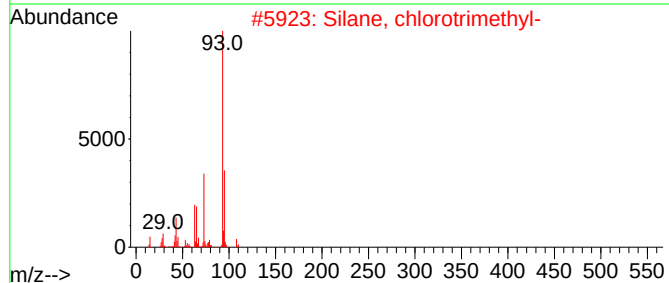
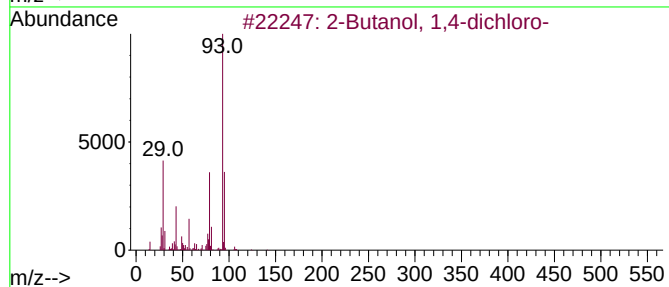
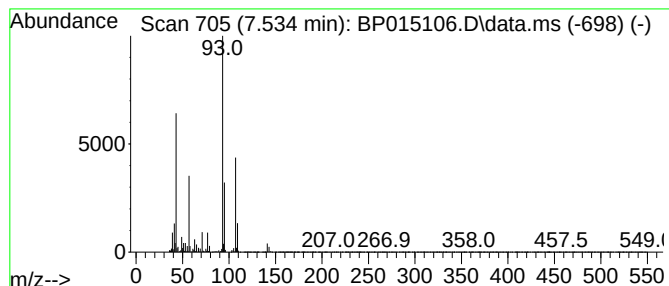
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.534	25.92 ng/ul	2573660	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 1,4-dichloro-			142	C4H8Cl2O	002419-74-1	38
2	Silane, chlorotrimethyl-			108	C3H9ClSi	000075-77-4	37
3	3-Pyrrolidinecarboxylic acid, 5-...			220	C11H12N2O3	1000337-74-3	9
4	Silane, chlorotrimethyl-			108	C3H9ClSi	000075-77-4	9
5	Silane, chlorotrimethyl-			108	C3H9ClSi	000075-77-4	8



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

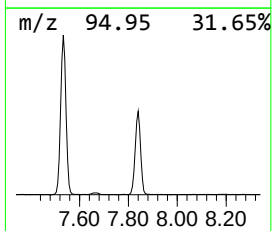
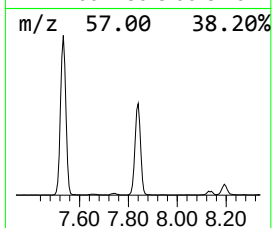
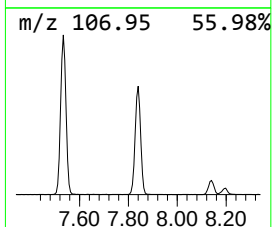
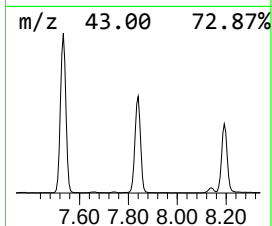
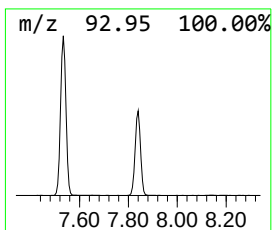
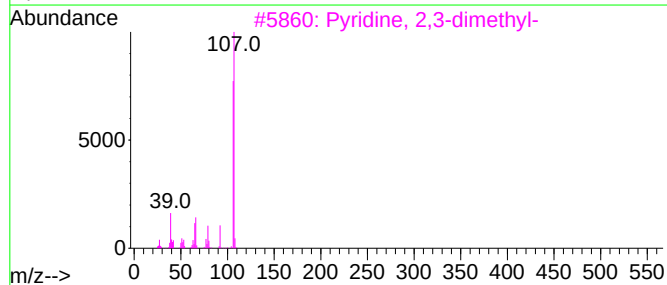
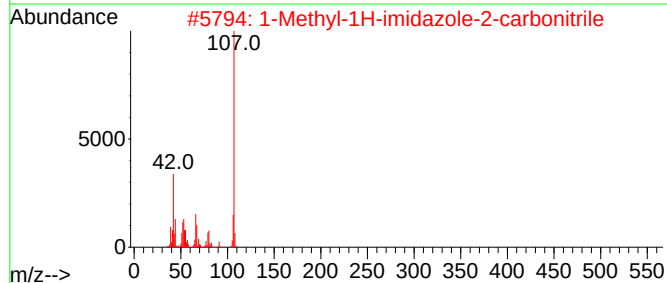
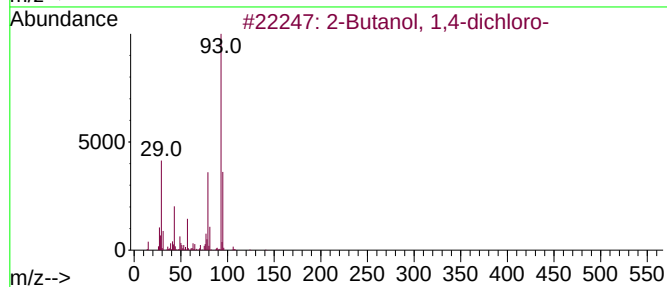
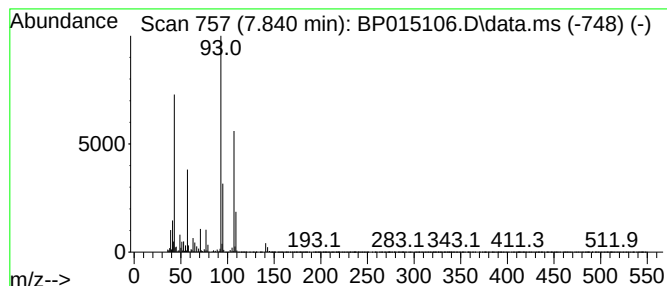
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.840	15.51 ng/ul	1540500	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 1,4-dichloro-			142	C4H8Cl2O	002419-74-1	40
2	1-Methyl-1H-imidazole-2-carbonit...			107	C5H5N3	1000306-60-1	9
3	Pyridine, 2,3-dimethyl-			107	C7H9N	000583-61-9	9
4	Silane, chlorotrimethyl-			108	C3H9ClSi	000075-77-4	9
5	Pyridine, 2,3-dimethyl-			107	C7H9N	000583-61-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP05106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

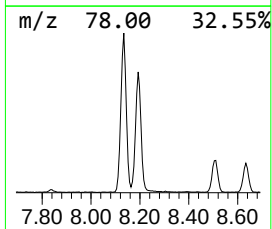
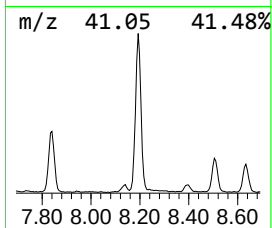
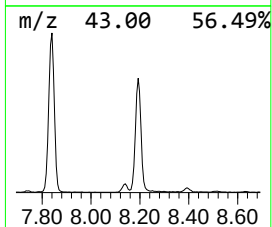
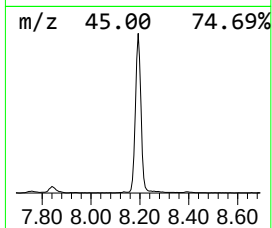
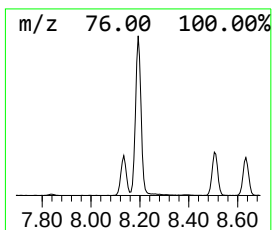
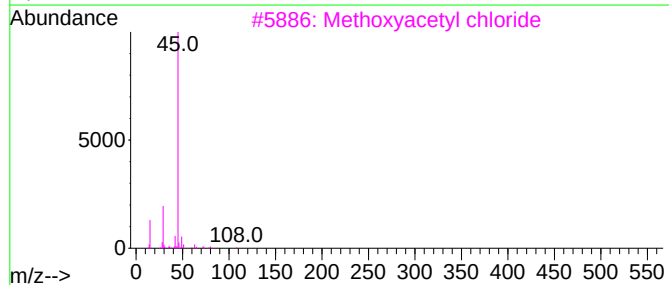
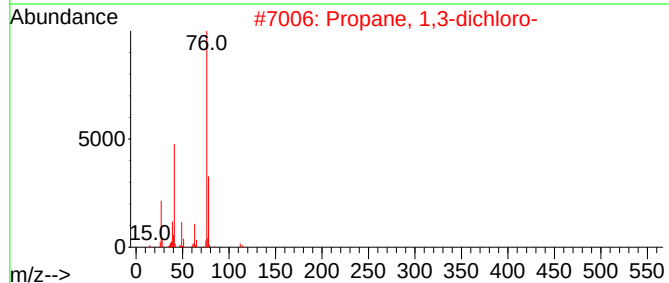
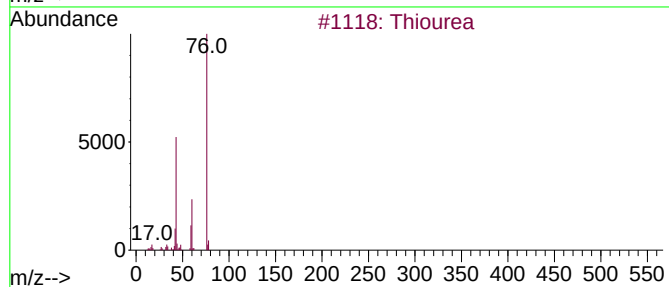
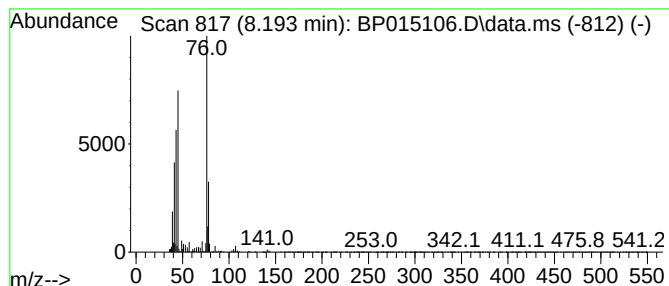
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.193	13.25 ng/ul	1315170	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiourea	76	CH4N2S	000062-56-6	42
2			Propane, 1,3-dichloro-	112	C3H6Cl2	000142-28-9	38
3			Methoxyacetyl chloride	108	C3H5ClO2	038870-89-2	10
4			Thiourea	76	CH4N2S	000062-56-6	9
5			Methoxyacetyl chloride	108	C3H5ClO2	038870-89-2	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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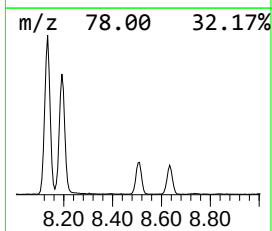
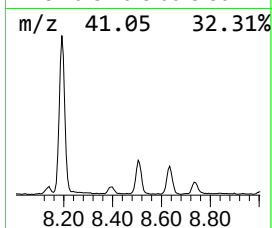
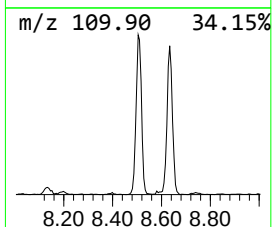
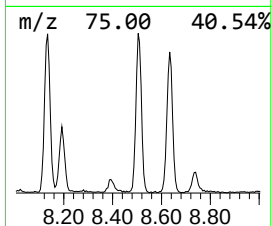
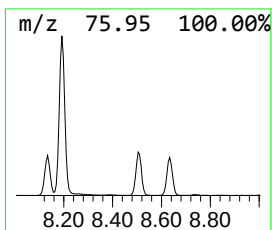
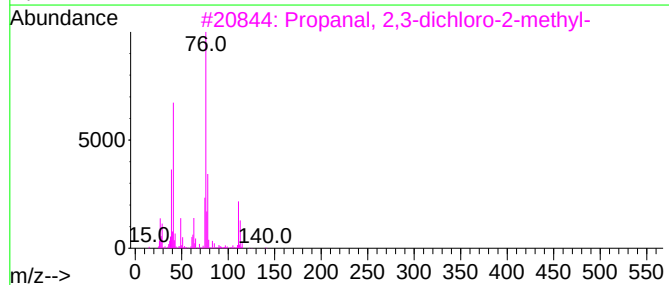
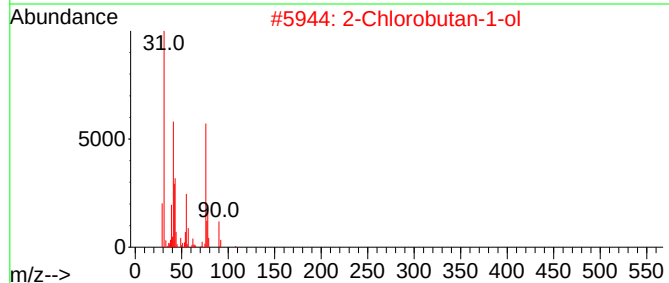
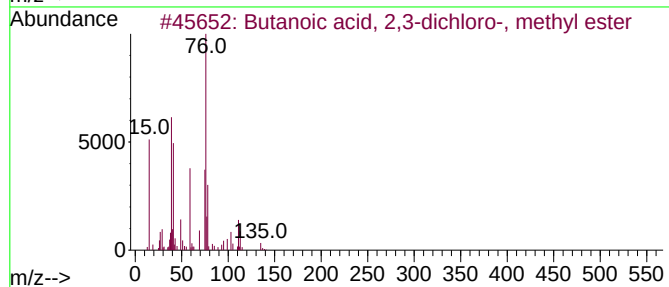
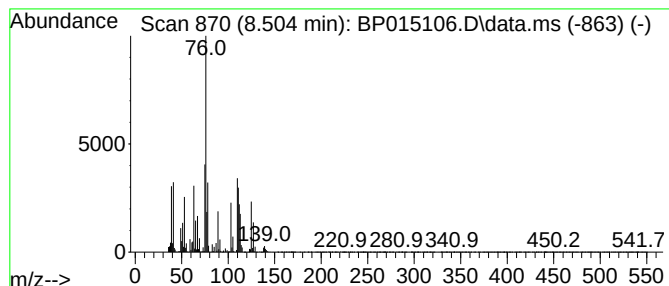
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Butanoic acid, 2,3-dichloro... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.504	5.94 ng/ul	589805	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 2,3-dichloro-, me...	170	C5H8Cl2O2	054460-97-8	56
2			2-Chlorobutan-1-ol	108	C4H9ClO	108269-46-1	40
3			Propanal, 2,3-dichloro-2-methyl-	140	C4H6Cl2O	010141-22-7	9
4			Propane, 1,3-dichloro-	112	C3H6Cl2	000142-28-9	9
5			Thiourea	76	CH4N2S	000062-56-6	5



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
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Sample : 02768-05
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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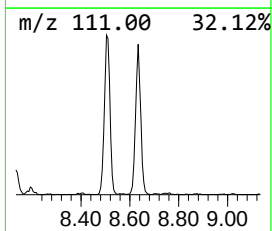
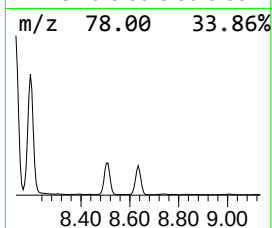
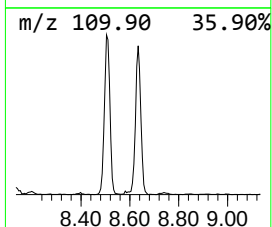
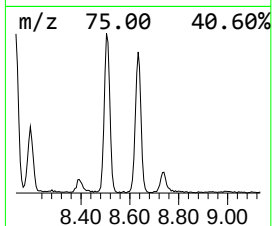
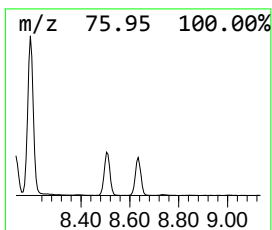
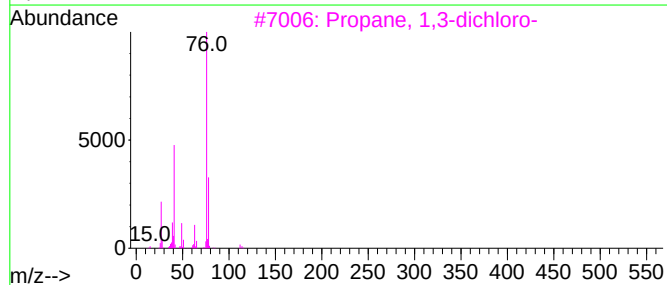
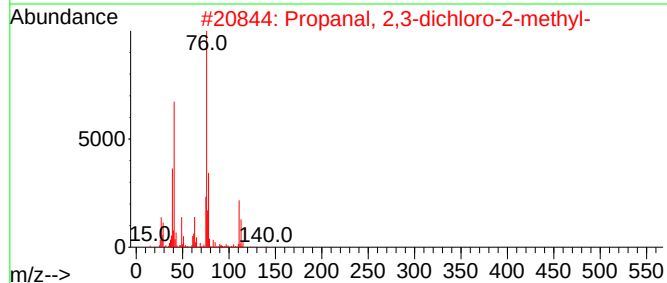
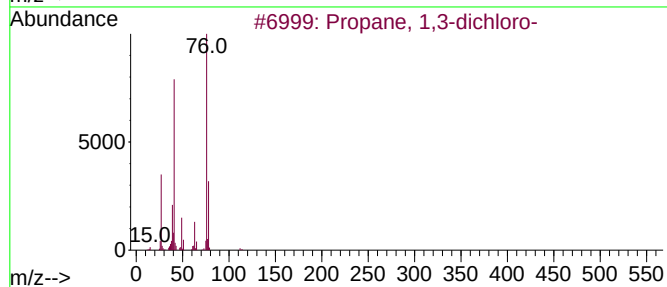
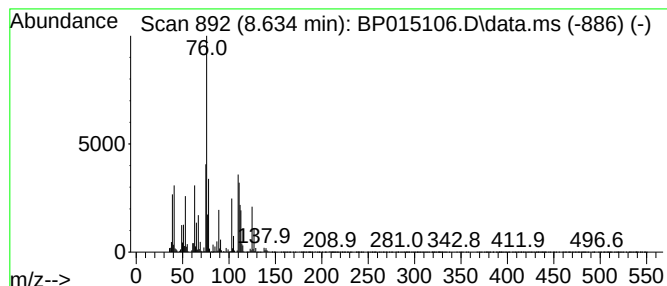
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Propane, 1,3-dichloro- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.634	4.82 ng/ul	479056	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,3-dichloro-	112	C3H6Cl2	000142-28-9	50
2			Propanal, 2,3-dichloro-2-methyl-	140	C4H6Cl2O	010141-22-7	45
3			Propane, 1,3-dichloro-	112	C3H6Cl2	000142-28-9	40
4			1H-Imidazole, 5-chloro-1-methyl-...	161	C4H4ClN3O2	004897-25-0	40
5			Butanoic acid, 2,3-dichloro-, me...	170	C5H8Cl2O2	054460-97-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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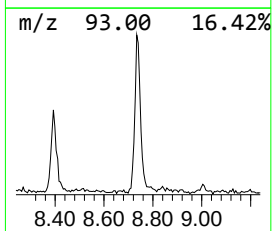
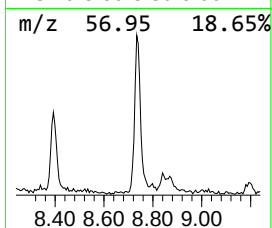
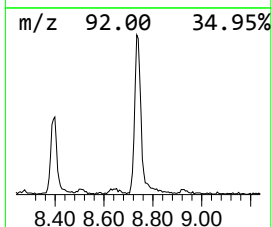
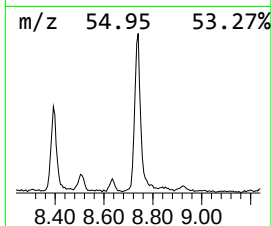
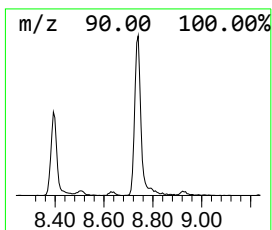
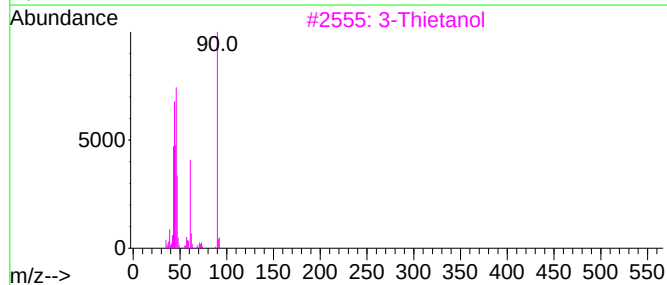
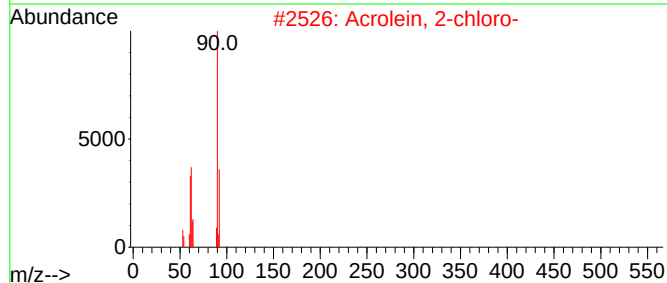
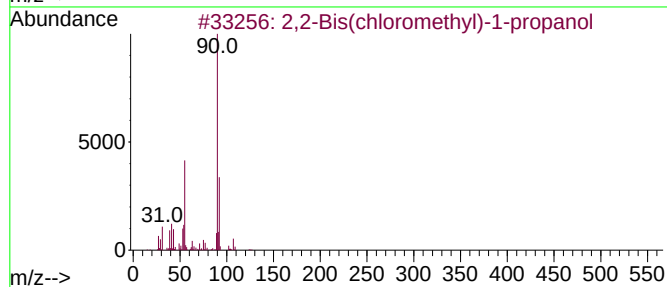
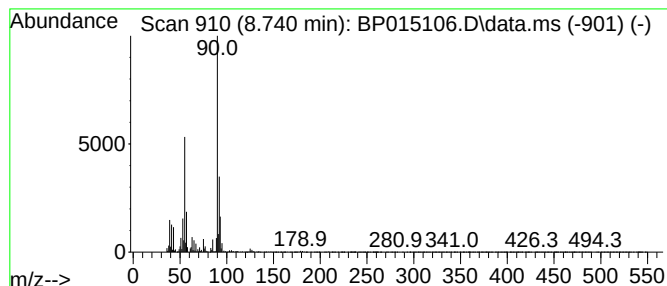
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 2,2-Bis(chloromethyl)-1-pro... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.740	3.41 ng/ul	338425	1,4-Dichlorobenzene-d4	8.134

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	64
2		Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	40
3		3-Thietanol	90	C3H6OS	010304-16-2	9
4		Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	9
5		Butanedioic acid, 2,3-dihydroxy-...	178	C6H10O6	000608-68-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
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Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CONN3

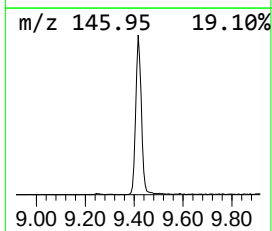
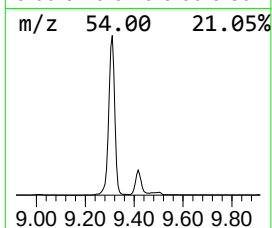
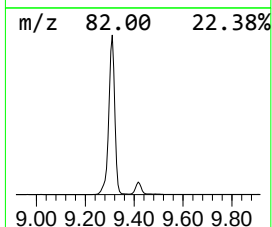
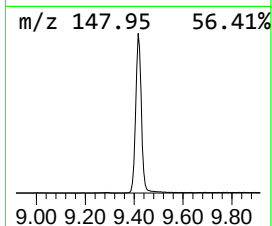
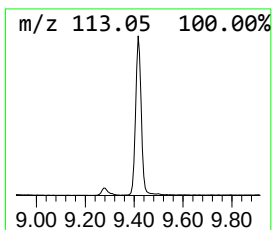
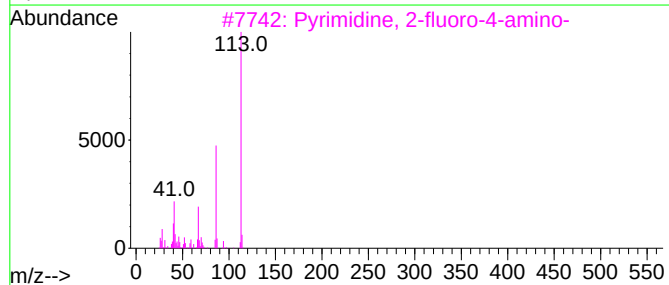
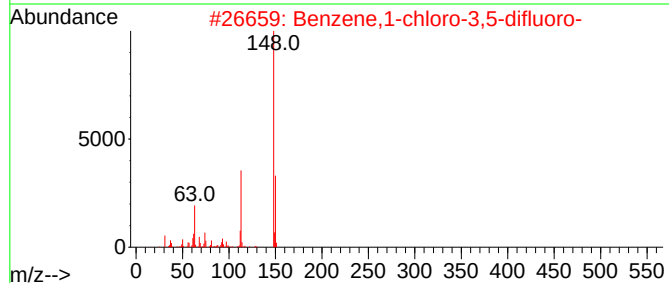
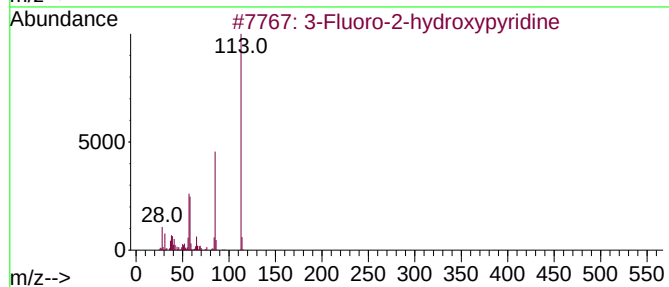
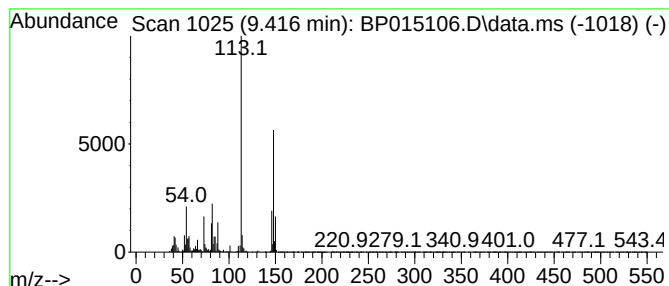
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-07 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.416	17.09 ng/ul	1696810	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Fluoro-2-hydroxypyridine	113	C5H4FN0	001547-29-1	25
2			Benzene,1-chloro-3,5-difluoro-	148	C6H3ClF2	001435-43-4	25
3			Pyrimidine, 2-fluoro-4-amino-	113	C4H4FN3	096548-91-3	22
4			Cyclopropanecarboxylic acid, 2,2...	154	C9H14O2	074685-76-0	17
5			1H-Pyrazole, 4-nitro-	113	C3H3N3O2	002075-46-9	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

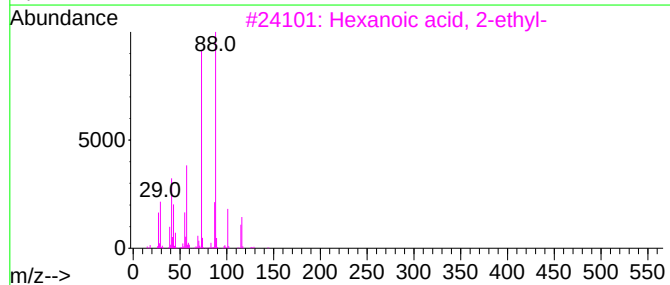
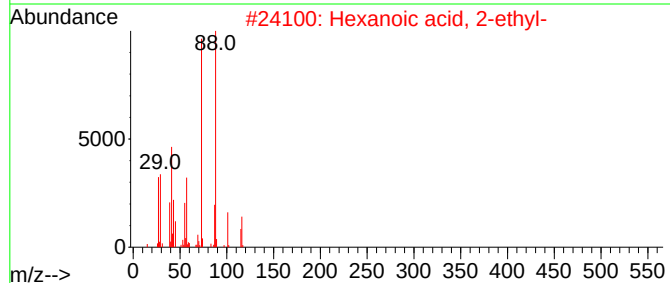
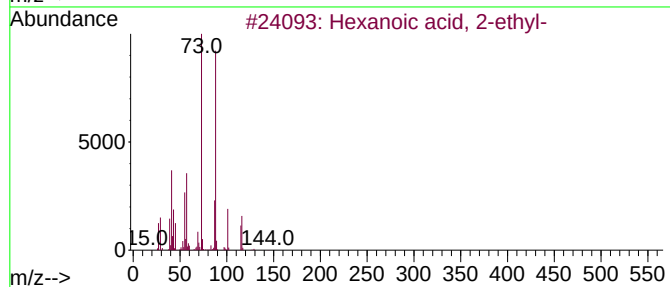
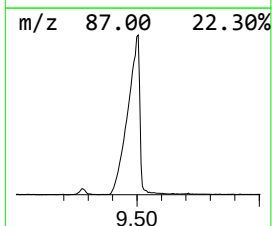
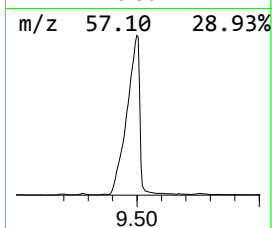
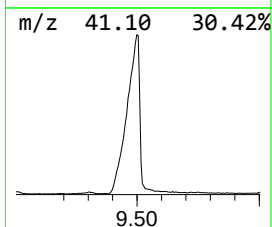
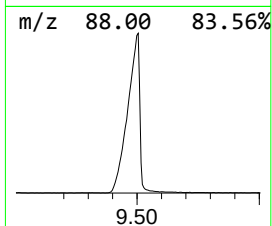
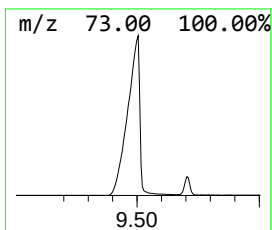
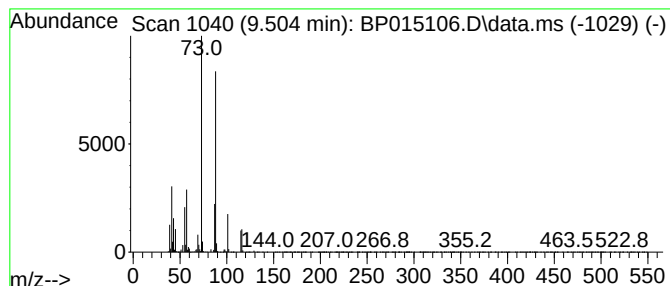
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Hexanoic acid, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.504	45.44 ng/ul	4512190	1,4-Dichlorobenzene-d4	8.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

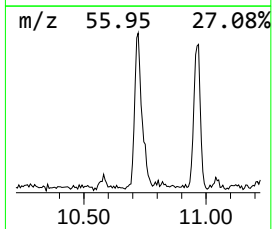
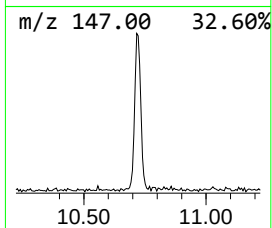
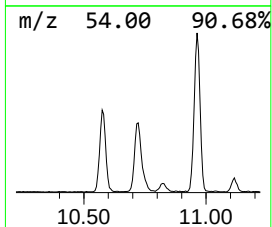
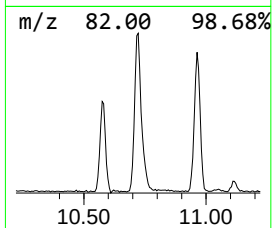
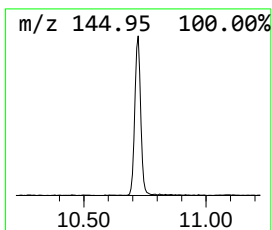
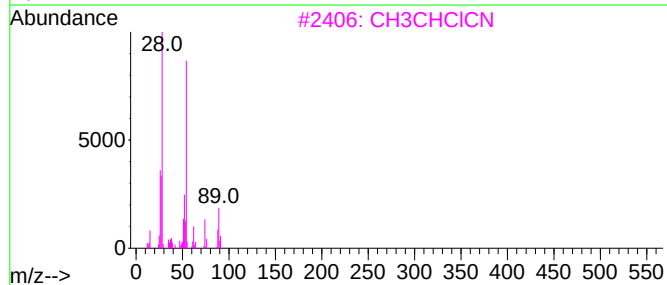
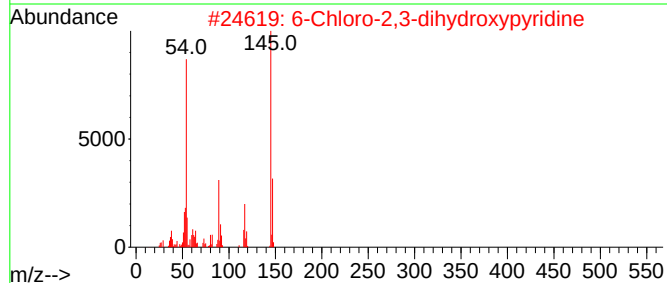
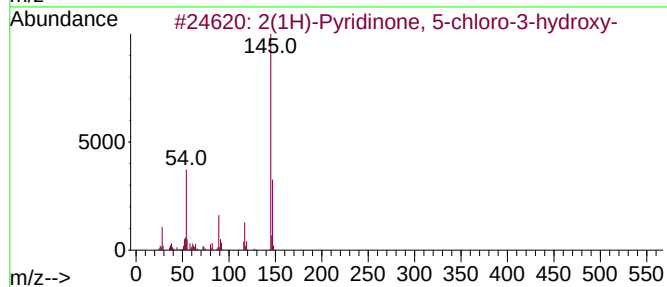
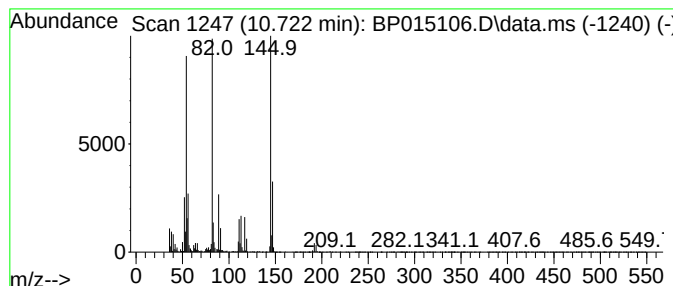
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 2(1H)-Pyridinone, 5-chloro-... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.722	3.71 ng/ul	540815	Naphthalene-d8	10.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Pyridinone, 5-chloro-3-hyd...	145	C5H4ClN02	053233-89-9	62
2	6-Chloro-2,3-dihydroxypyridine	145	C5H4ClN02	1000341-18-9	58
3	CH3CHClCN	89	C3H4ClN	001617-17-0	50
4	Propanoic acid, 3-cyano-, methyl...	113	C5H7N02	004107-62-4	40
5	CH3CHClCN	89	C3H4ClN	001617-17-0	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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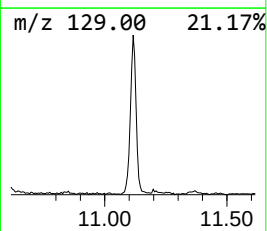
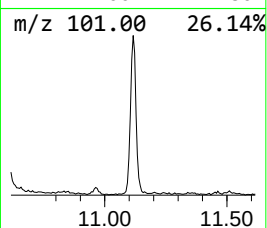
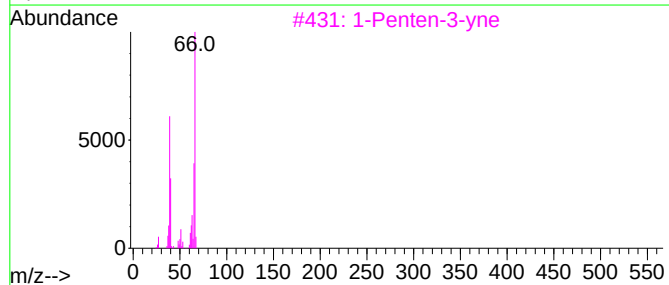
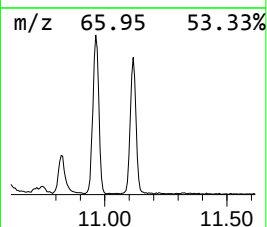
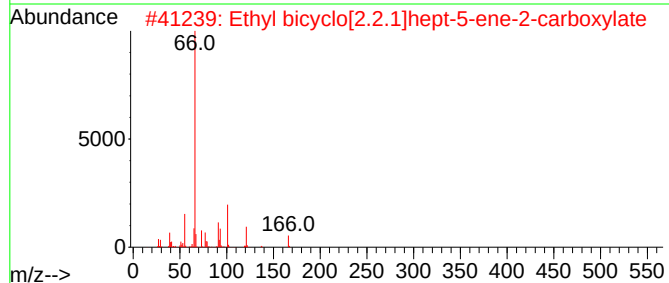
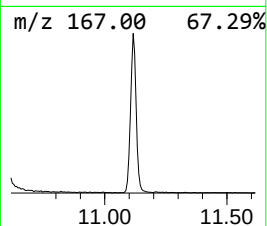
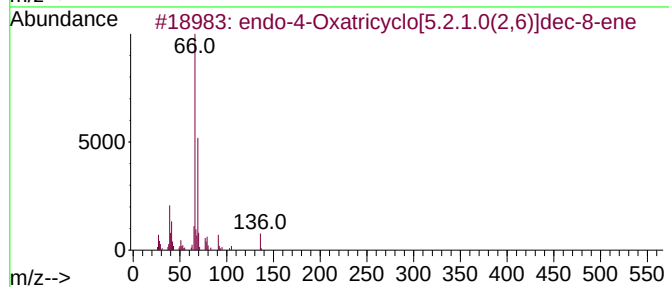
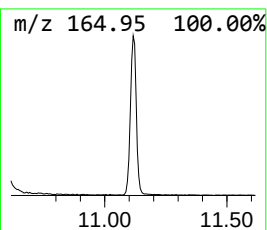
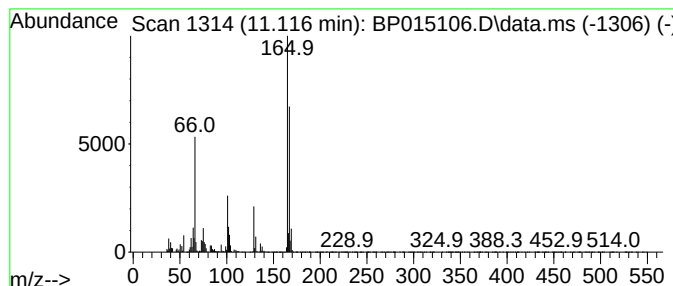
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 unknown-08 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.116	4.19 ng/ul	611175	Naphthalene-d8	10.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	endo-4-Oxatricyclo[5.2.1.0(2,6)]...	136	C9H12O	001528-23-0	35		
2	Ethyl bicyclo[2.2.1]hept-5-ene-2...	166	C10H14O2	010138-32-6	35		
3	1-Penten-3-yne	66	C5H6	000646-05-9	35		
4	2,6-Dichloro-3-fluoropyridine	165	C5H2Cl2FN	052208-50-1	30		
5	Bicyclo[2.2.1]hept-5-ene-2-carbo...	138	C8H10O2	000120-74-1	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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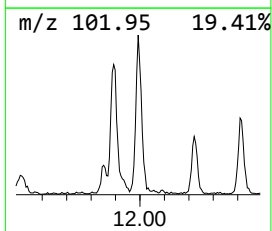
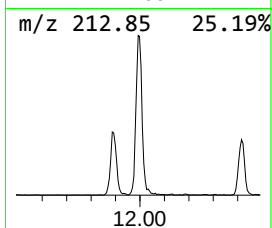
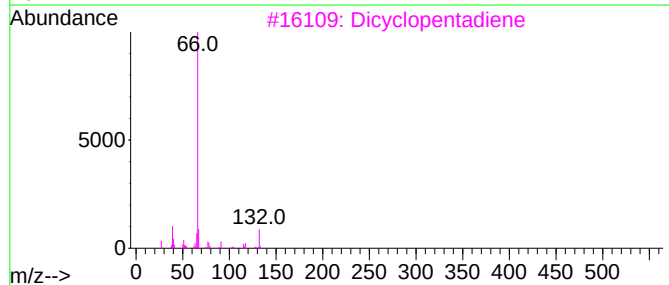
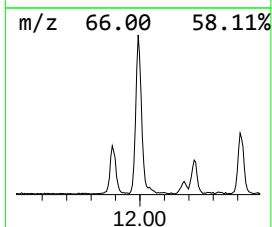
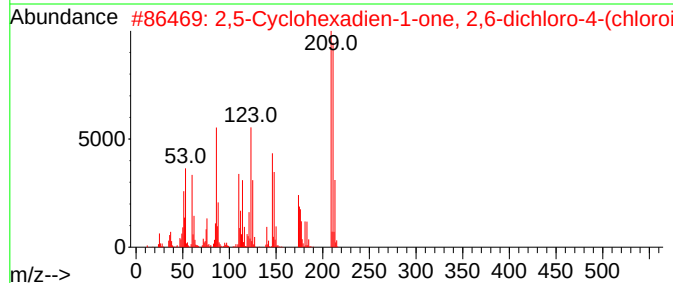
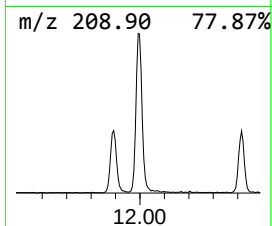
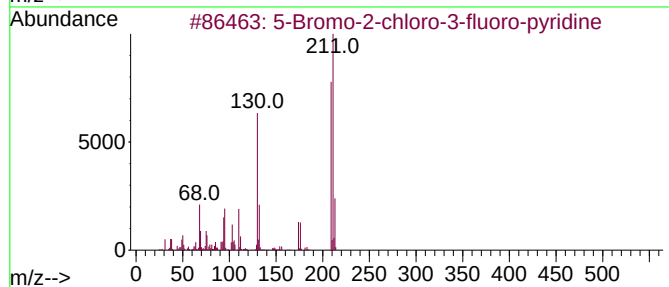
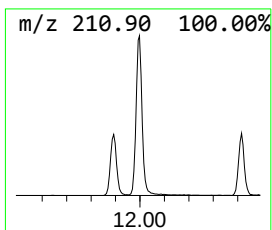
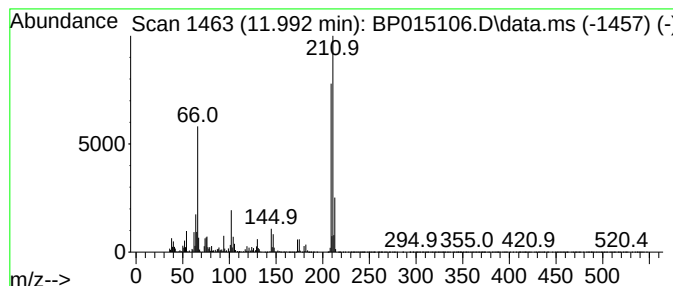
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 unknown-09 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.992	3.84 ng/ul	560873	Naphthalene-d8	10.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Bromo-2-chloro-3-fluoro-pyridine	209	C5H2BrClFN	831203-13-5	38
2		2,5-Cyclohexadien-1-one, 2,6-dic...	209	C6H2Cl3NO	000101-38-2	35
3		Dicyclopentadiene	132	C10H12	000077-73-6	25
4		5-Chlorobicyclo[2.2.1]hept-2-ene...	128	C7H9Cl	1000436-27-6	22
5		2-Norbornene	94	C7H10	000498-66-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

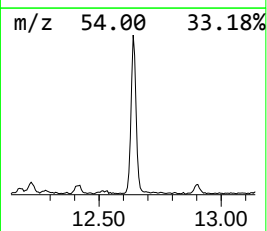
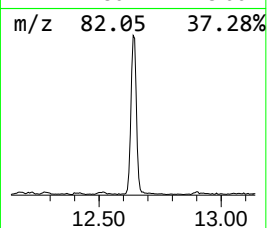
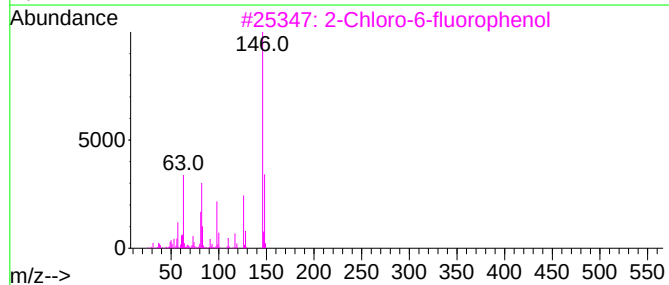
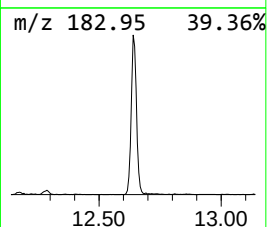
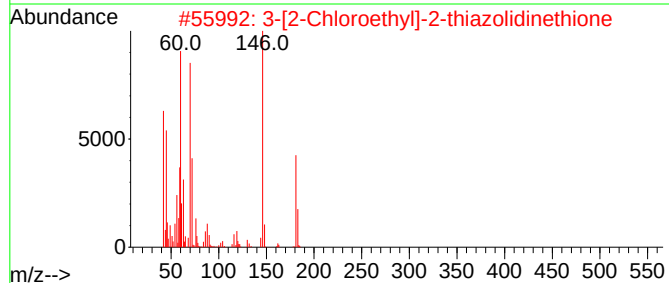
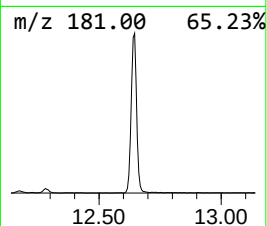
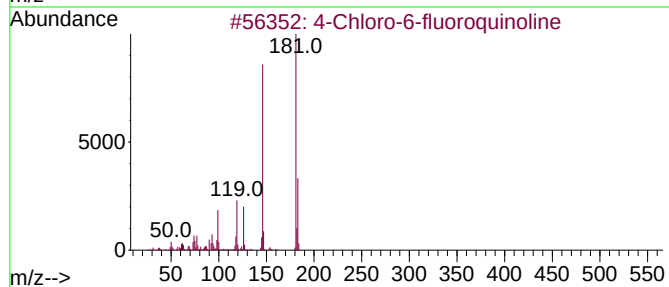
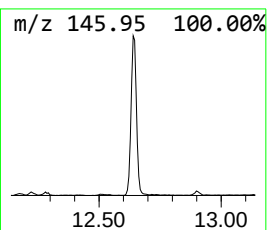
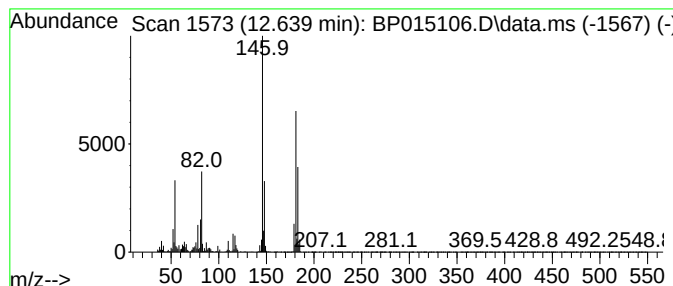
Instrument :
BNA_P
ClientSampleId :
CONN3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-10 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.639	6.79 ng/ul	991092	Naphthalene-d8	10.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-6-fluoroquinoline	181	C9H5ClFN	000391-77-5	43
2	3-[2-Chloroethyl]-2-thiazolidine...	181	C5H8ClNS2	089416-25-1	40
3	2-Chloro-6-fluorophenol	146	C6H4ClFO	002040-90-6	38
4	2-Chloro-4-fluorophenol	146	C6H4ClFO	001996-41-4	32
5	2-Chloro-5-trifluoromethylpyridine	181	C6H3ClF3N	052334-81-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

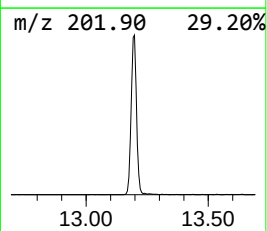
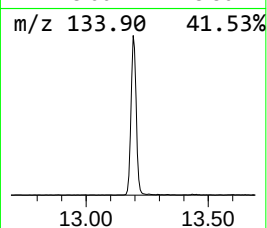
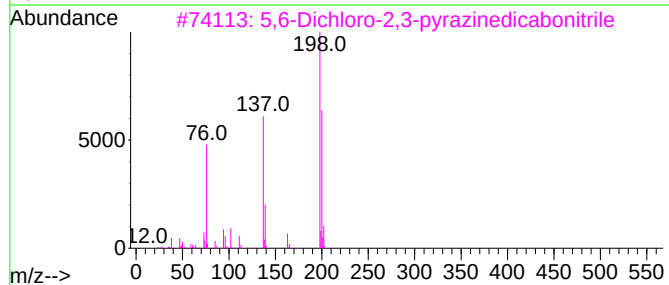
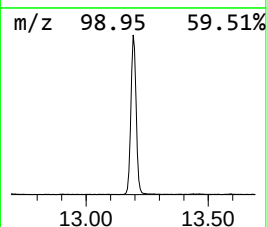
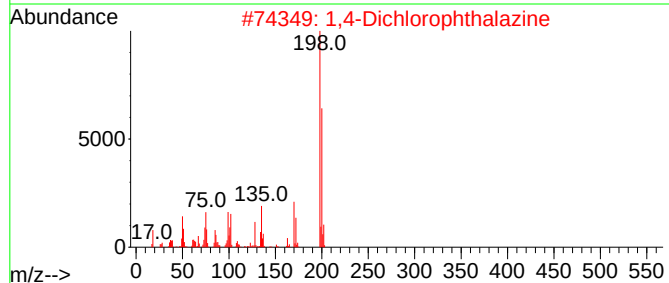
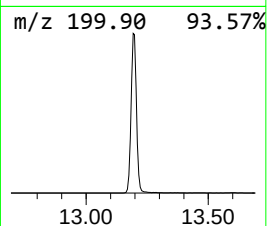
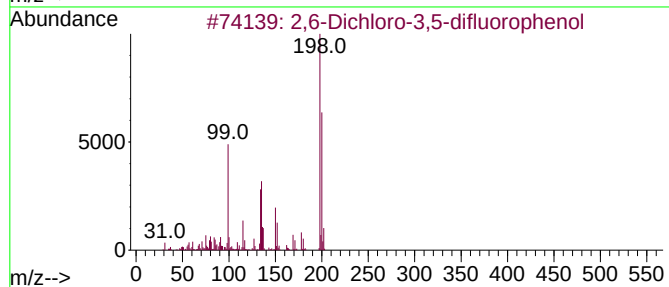
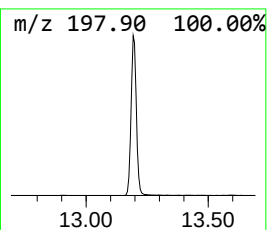
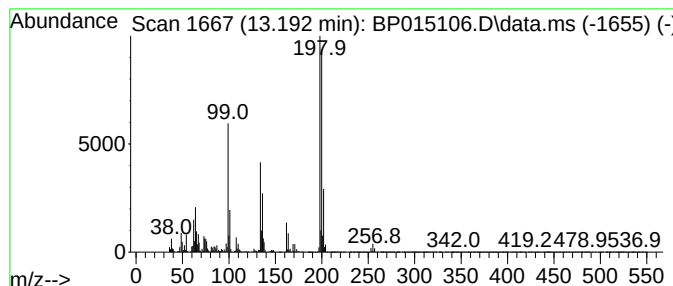
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 2,6-Dichloro-3,5-difluoroph... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.192	8.29 ng/ul	1710140	Acenaphthene-d10			14.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Dichloro-3,5-difluorophenol	198	C6H2Cl2F2O	063418-08-6	58	
2	1,4-Dichlorophthalazine	198	C8H4Cl2N2	004752-10-7	46	
3	5,6-Dichloro-2,3-pyrazinedicabon...	198	C6Cl2N4	056413-95-7	35	
4	1-(3-Bromophenyl)piperazine	240	C10H13BrN2	031197-30-5	27	
5	1,5-Naphthyridine, 4,8-dichloro-	198	C8H4Cl2N2	028252-80-4	22	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

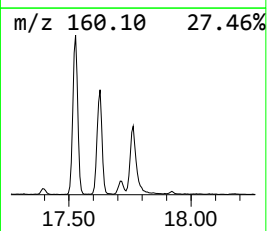
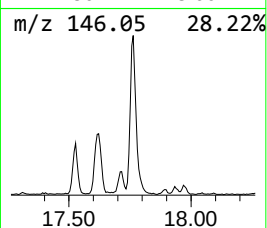
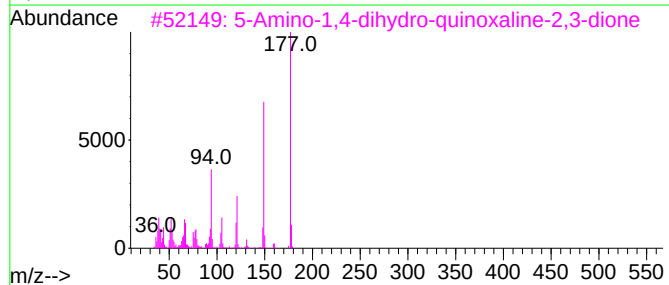
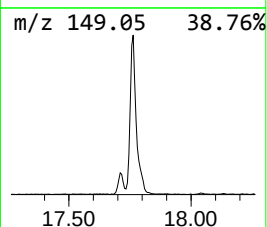
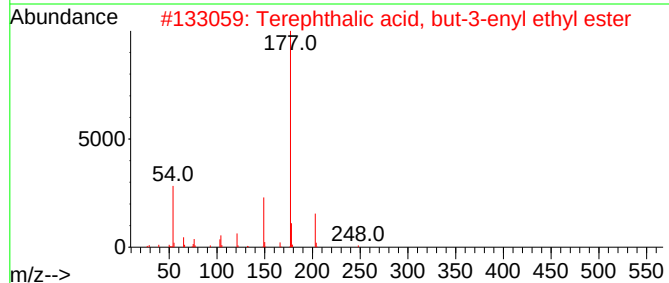
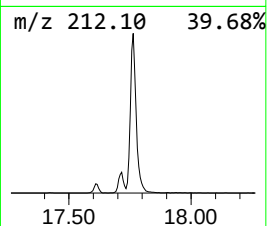
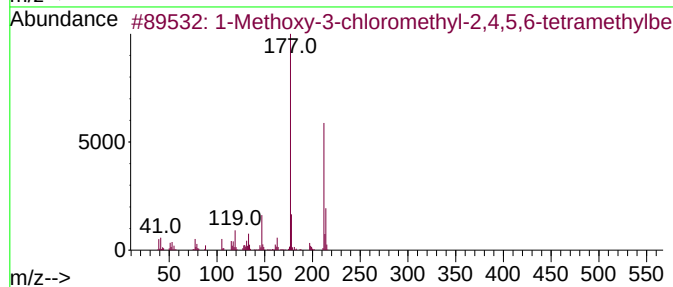
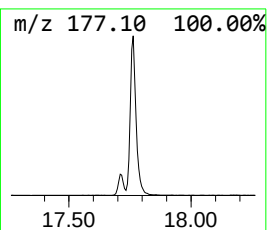
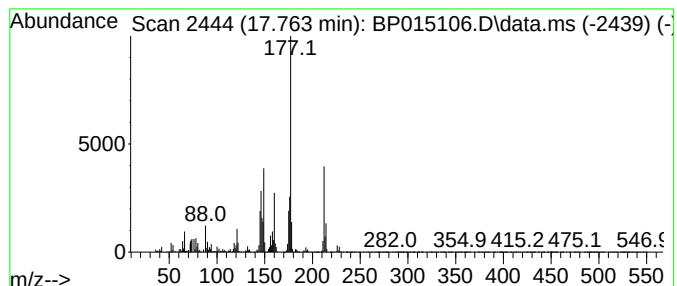
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 unknown-11 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.763	8.79 ng/ul	2182700	Phenanthrene-d10	17.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methoxy-3-chloromethyl-2,4,5,6...	212	C12H17ClO	073942-80-0	43
2			Terephthalic acid, but-3-enyl et...	248	C14H16O4	1000356-33-3	35
3			5-Amino-1,4-dihydro-quinoxaline-...	177	C8H7N3O2	076097-87-5	35
4			Terephthalic acid, ethyl phenyl ...	270	C16H14O4	1000324-05-8	27
5			Spiro[cyclopenta[d]-1,3,2-dioxab...	206	C12H19B02	062238-26-0	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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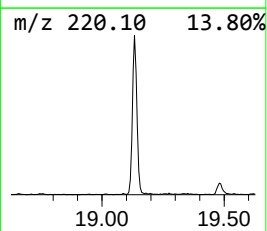
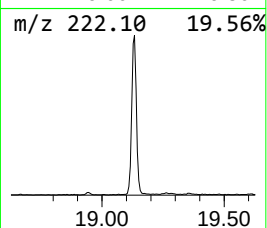
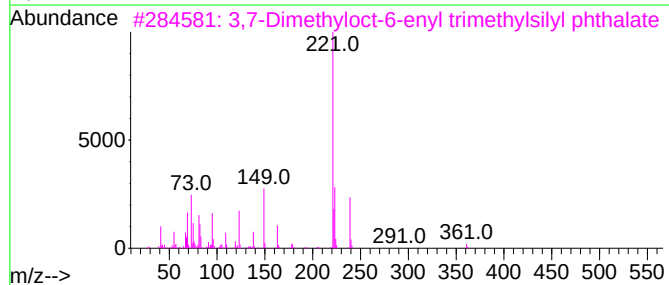
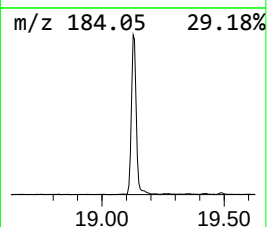
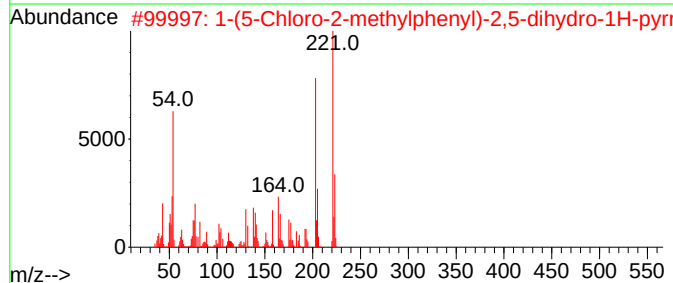
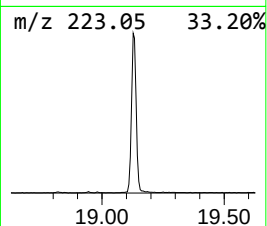
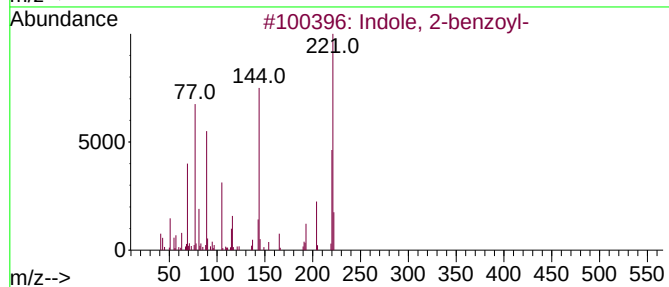
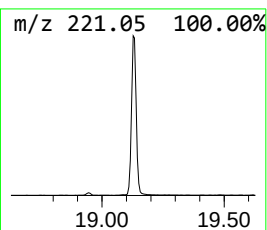
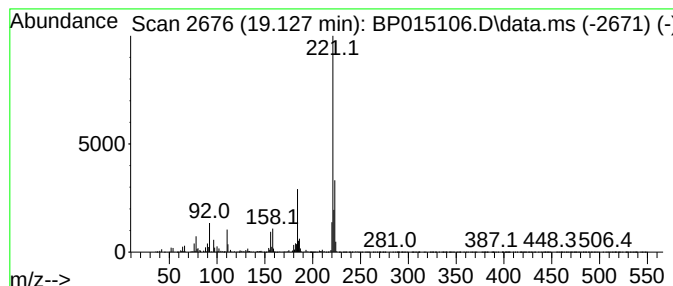
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 unknown-12 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.127	5.74 ng/ul	1423840	Phenanthrene-d10	17.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indole, 2-benzoyl-	221	C15H11NO	001022-86-2	43
2			1-(5-Chloro-2-methylphenyl)-2,5-...	221	C11H8ClNO2	058670-26-1	40
3			3,7-Dimethyloct-6-enyl trimethyl...	376	C21H32O4Si	1000373-64-6	40
4			4-Chloro-3,5-difluorophenol, TMS	236	C9H11ClF2OSi	1000508-21-4	33
5			2-Methyl-thieno[2',3':3,4]benzo[...	221	C10H7NOS2	294668-48-7	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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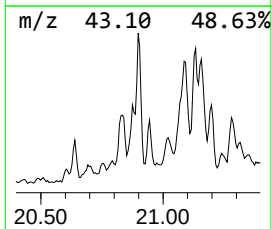
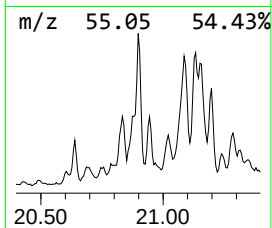
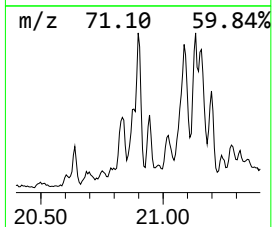
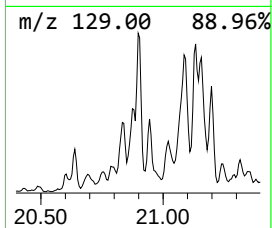
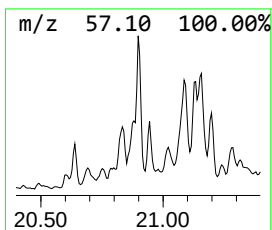
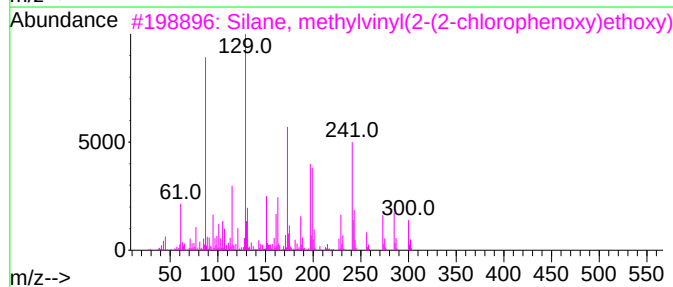
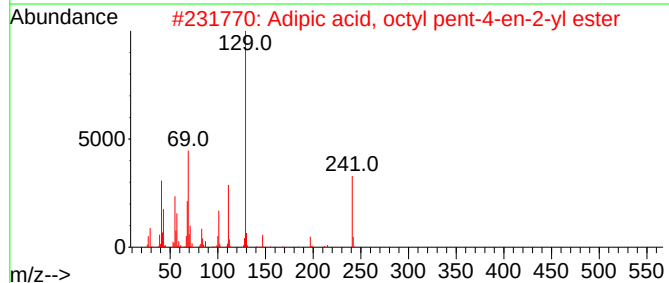
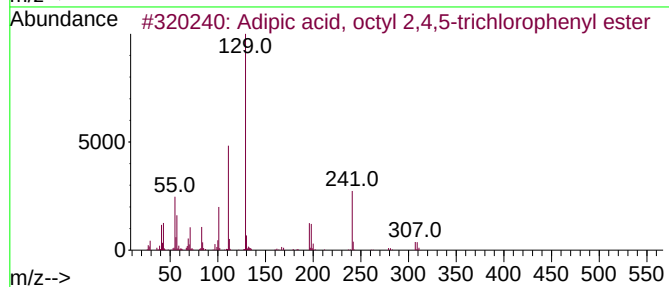
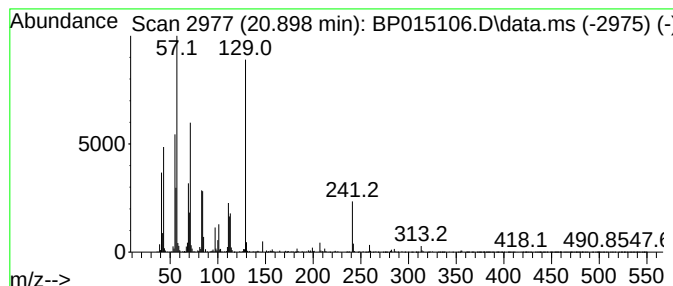
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Adipic acid, octyl 2,4,5-tr... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.898	3.74 ng/ul	944295	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adipic acid, octyl 2,4,5-trichlo...	436	C20H27Cl3O4	1000353-86-1	50		
2	Adipic acid, octyl pent-4-en-2-y...	326	C19H34O4	1000354-12-4	49		
3	Silane, methylvinyl(2-(2-chlorop...	300	C14H21ClO3Si	1000420-77-6	30		
4	Adipic acid, octyl 2,3,5-trichlo...	436	C20H27Cl3O4	1000353-87-3	27		
5	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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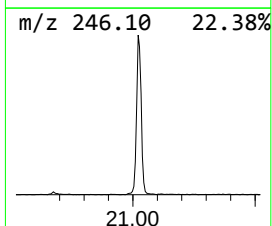
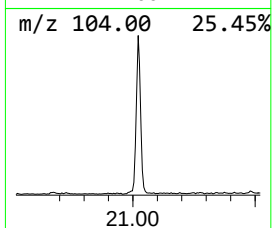
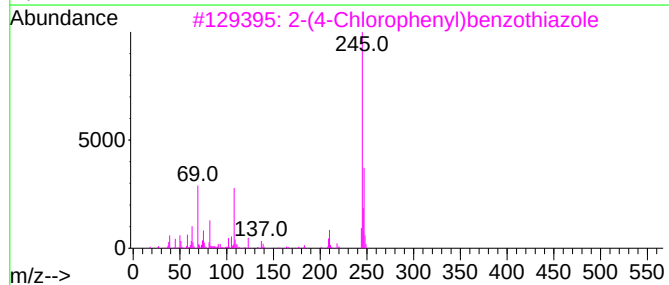
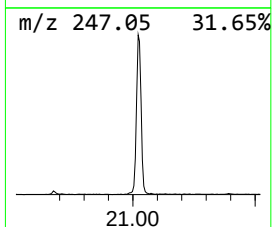
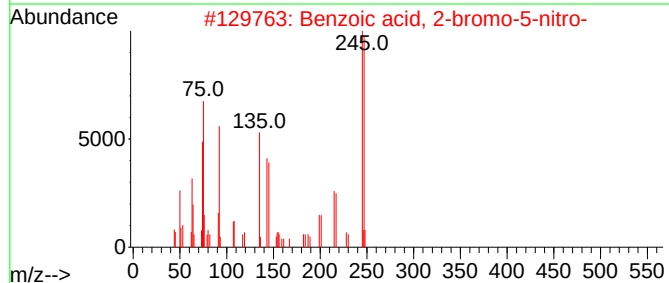
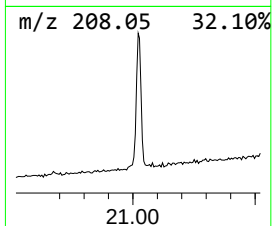
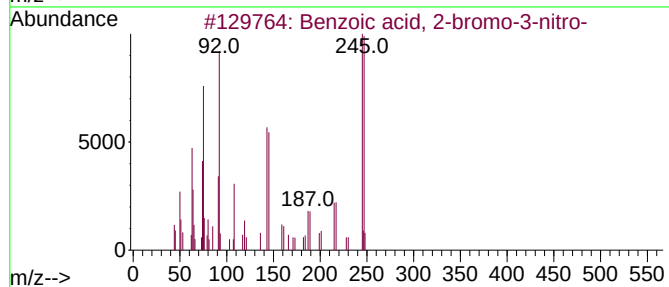
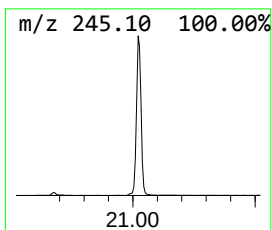
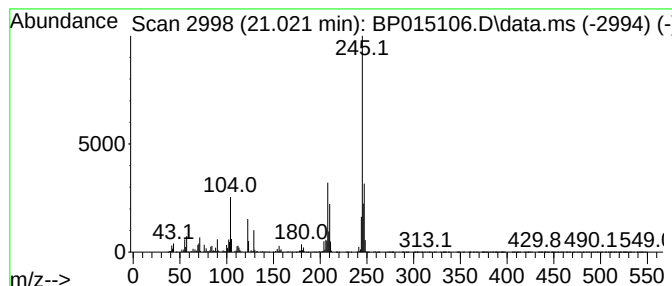
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Benzoic acid, 2-bromo-3-nitro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.021	5.57 ng/ul	1406510	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2-bromo-3-nitro-	245	C7H4BrNO4	000573-54-6	89
2			Benzoic acid, 2-bromo-5-nitro-	245	C7H4BrNO4	000943-14-6	89
3			2-(4-Chlorophenyl)benzothiazole	245	C13H8ClNS	006265-91-4	62
4			2-(4-chlorophenyl)-5-fluoro-1H-i...	245	C14H9ClFN	881040-32-0	58
5			4-Chloro-2-methyl-7-(trifluorome...	245	C11H7ClF3N	018529-09-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

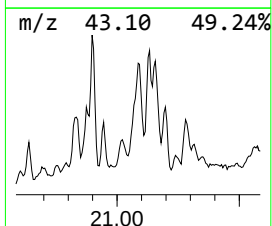
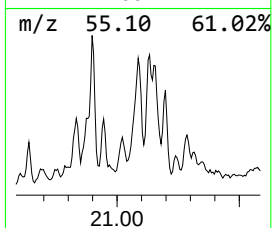
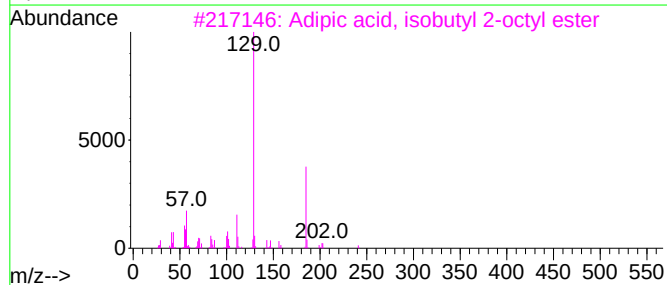
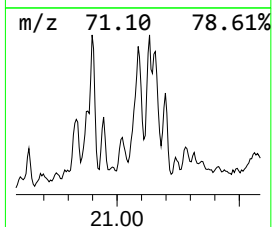
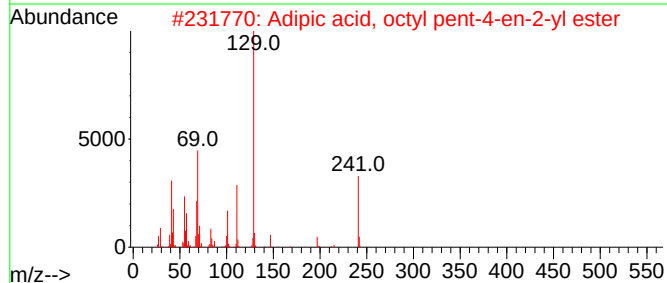
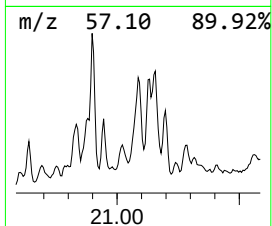
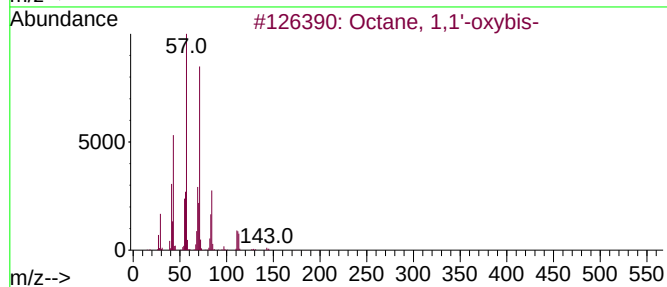
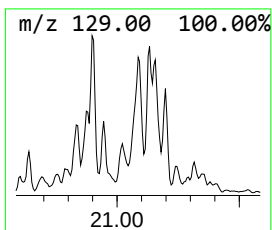
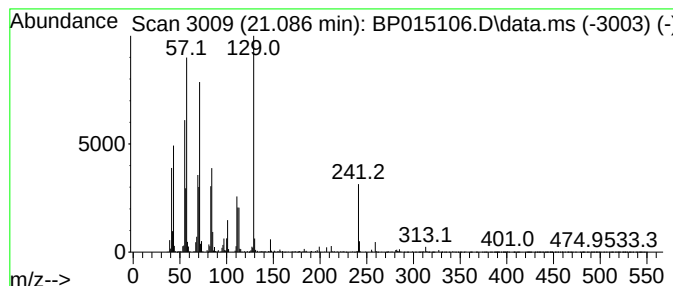
Instrument :
BNA_P
ClientSampleId :
CONN3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 35 unknown-13 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.086	4.45 ng/ul	1125250	Chrysene-d12	21.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	43
2	Adipic acid, octyl pent-4-en-2-yl ester	326	C19H34O4	1000354-12-4	38
3	Adipic acid, isobutyl 2-octyl ester	314	C18H34O4	1000324-44-5	27
4	3-Methyl-2-butenic acid, 2-tetradecyl ester	184	C10H16O3	1010279-23-1	27
5	Oxalic acid, 2-ethylhexyl isohexadecyl ester	286	C16H30O4	1000309-38-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP05106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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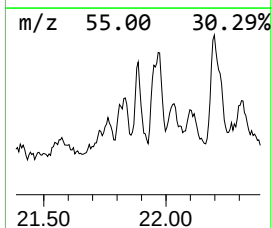
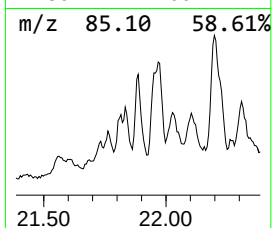
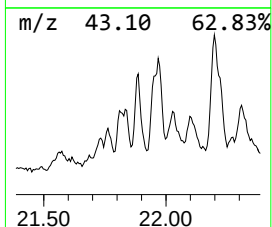
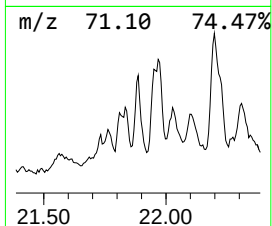
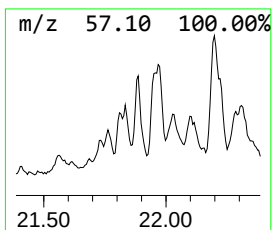
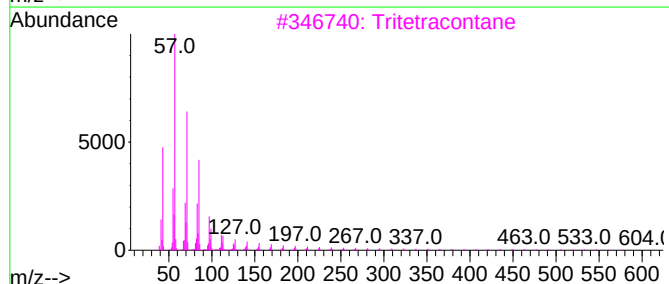
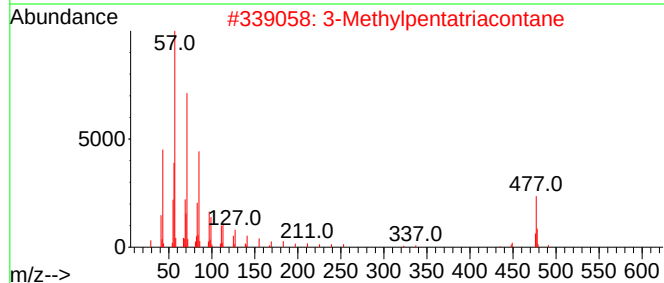
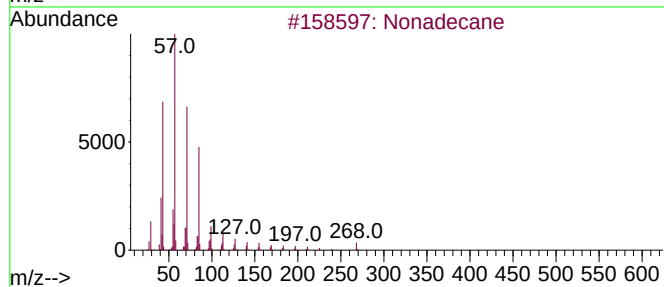
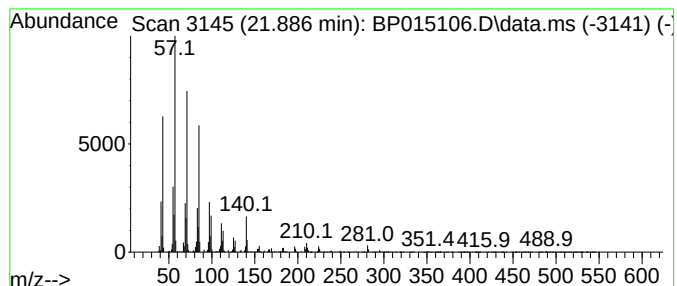
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.886	2.27 ng/ul	573430	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			3-Methylpentatriacontane	507	C36H74	078692-70-3	87
3			Tritetracontane	605	C43H88	007098-21-7	83
4			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	83
5			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

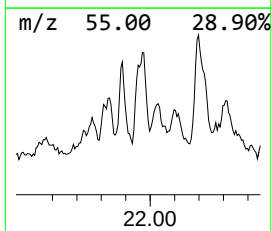
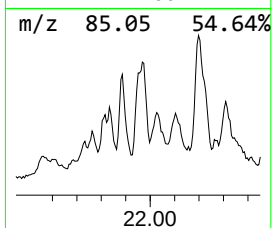
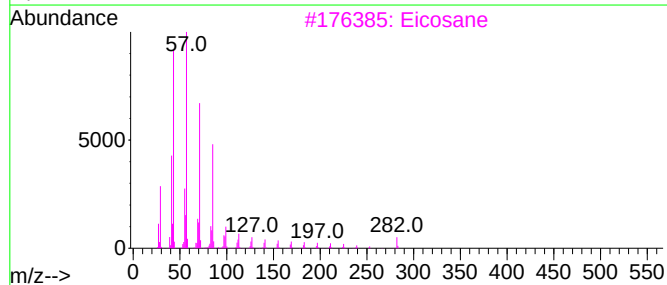
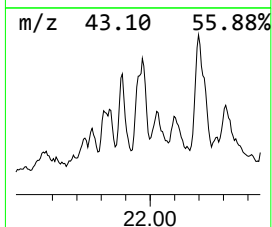
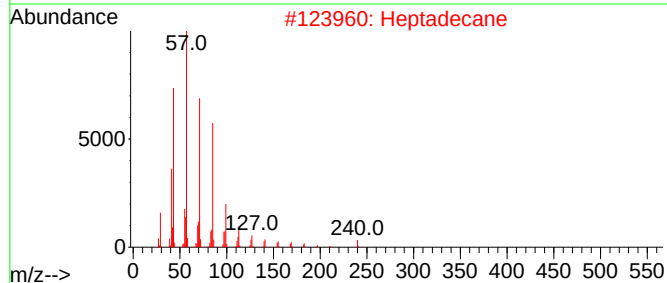
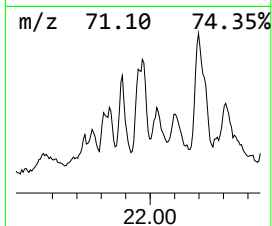
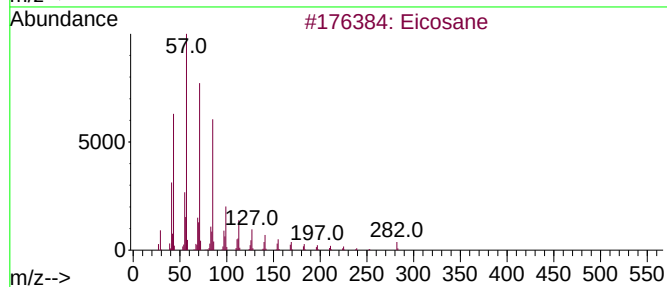
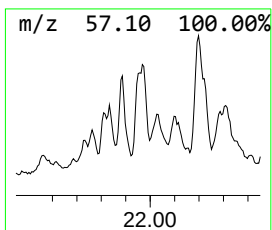
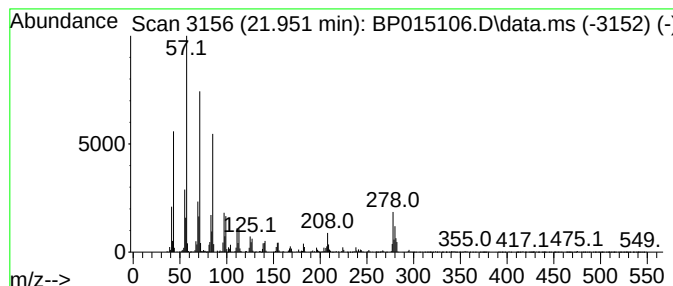
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.951	2.66 ng/ul	672100	Chrysene-d12		21.604	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	Heptadecane		240	C17H36	000629-78-7	93
3	Eicosane		282	C20H42	000112-95-8	89
4	Tetracosane		338	C24H50	000646-31-1	76
5	10-Methylnonadecane		282	C20H42	056862-62-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
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Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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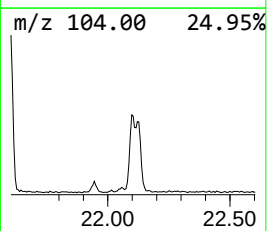
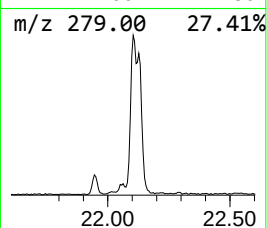
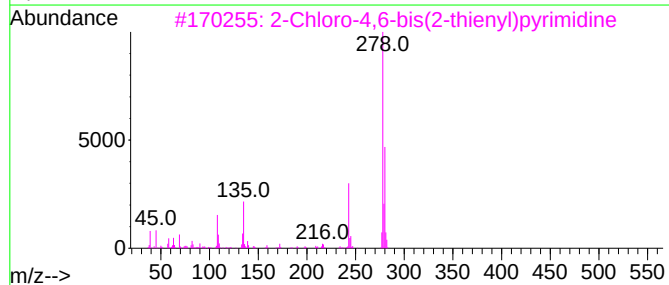
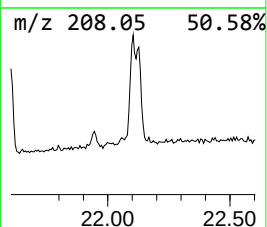
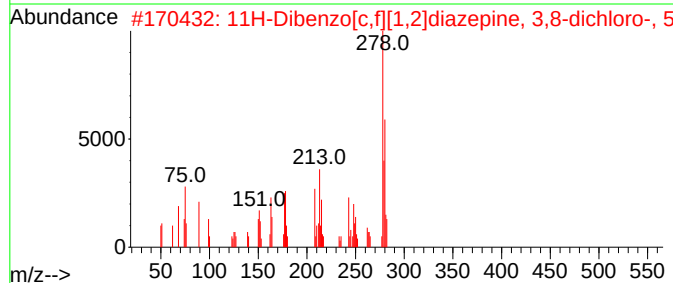
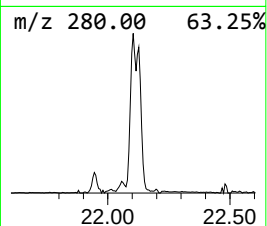
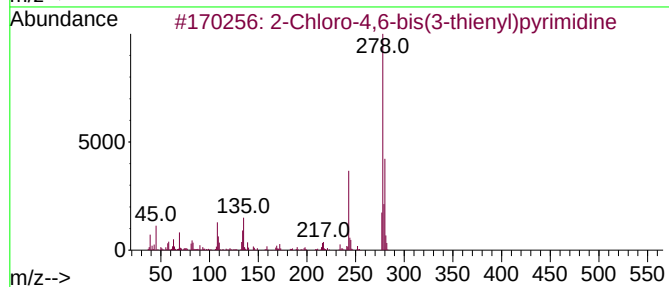
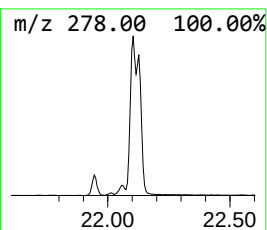
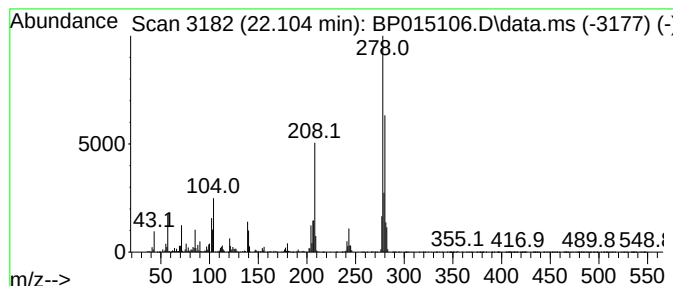
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 2-Chloro-4,6-bis(3-thienyl)... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.104	4.69 ng/ul	1185930	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-4,6-bis(3-thienyl)pyrim...	278	C12H7ClN2S2	131022-83-8	62
2			11H-Dibenzo[c,f][1,2]diazepine, ...	278	C13H8ClN2O	000958-71-4	53
3			2-Chloro-4,6-bis(2-thienyl)pyrim...	278	C12H7ClN2S2	124959-36-0	50
4			4-Bromophenoxathiin	278	C12H7BrOS	105906-37-4	43
5			Benzo[b]triphenylene	278	C22H14	000215-58-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

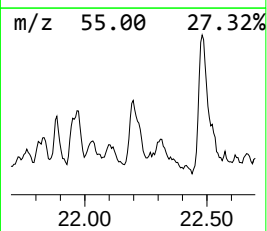
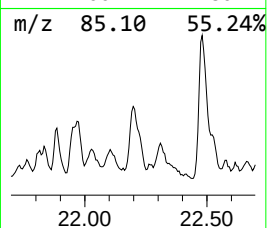
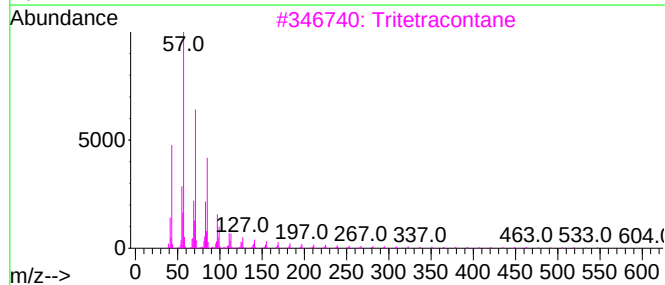
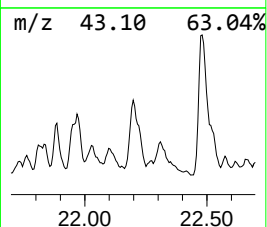
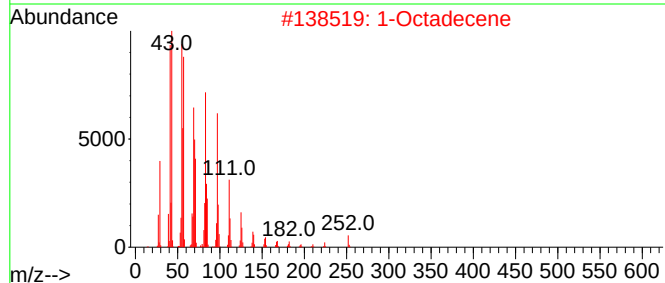
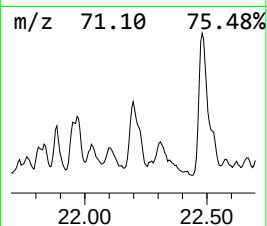
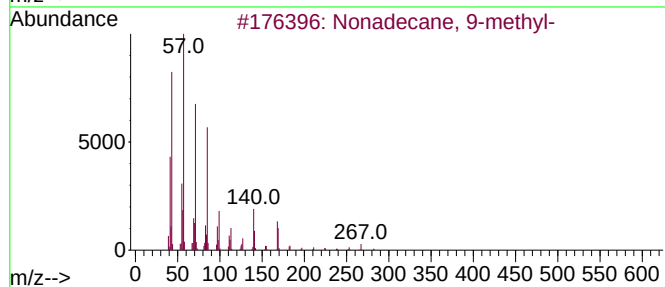
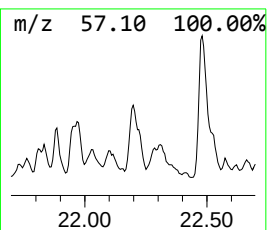
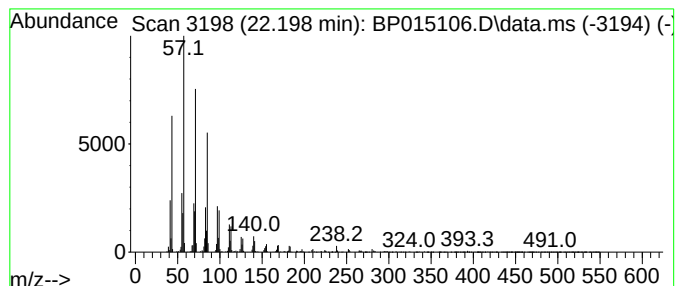
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.198	4.90 ng/ul	1238640	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
2			1-Octadecene	252	C18H36	000112-88-9	92
3			Tritetracontane	605	C43H88	007098-21-7	91
4			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	87
5			Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

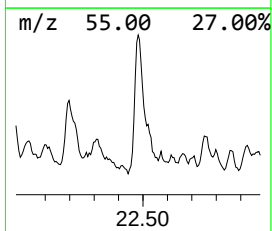
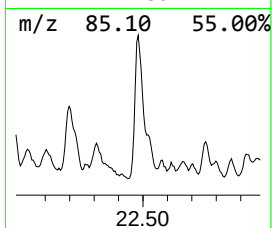
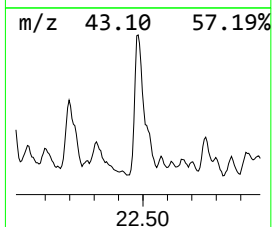
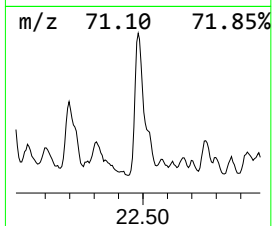
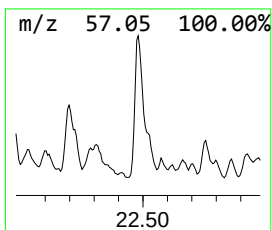
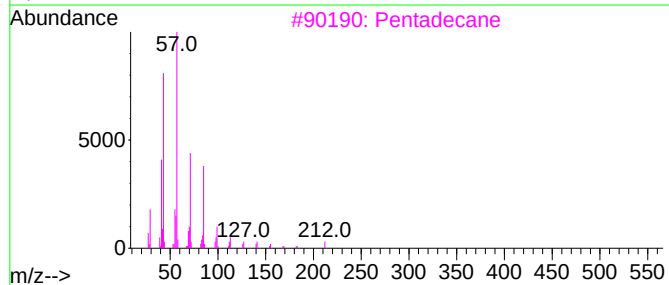
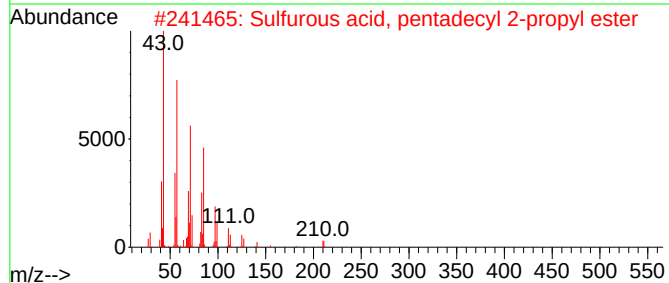
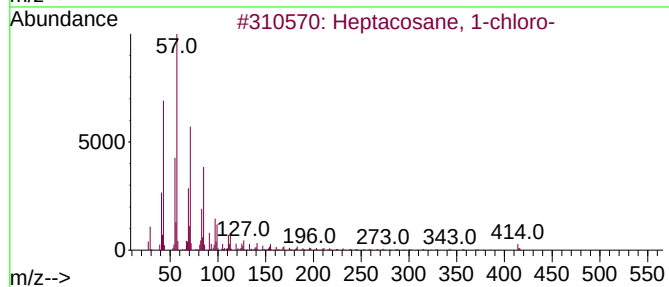
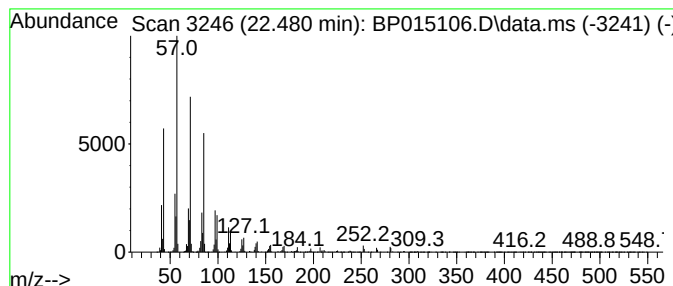
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Heptacosane, 1-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.480	11.70 ng/ul	2955400	Chrysene-d12	21.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
2			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	91
3			Pentadecane	212	C15H32	000629-62-9	89
4			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	83
5			Hexacosane	366	C26H54	000630-01-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP05106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

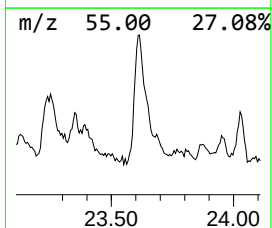
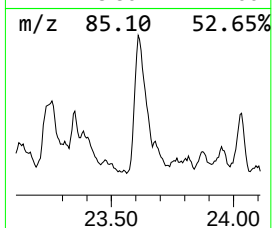
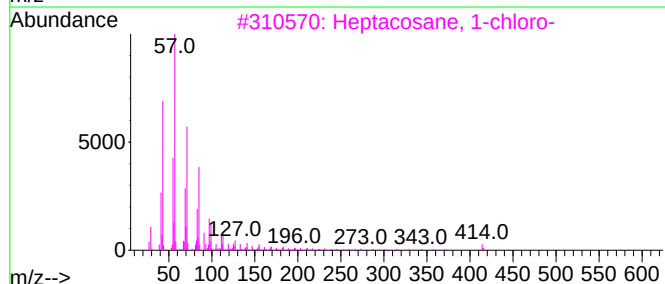
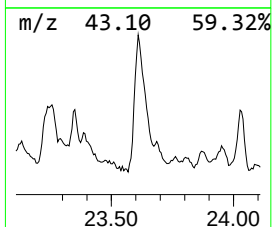
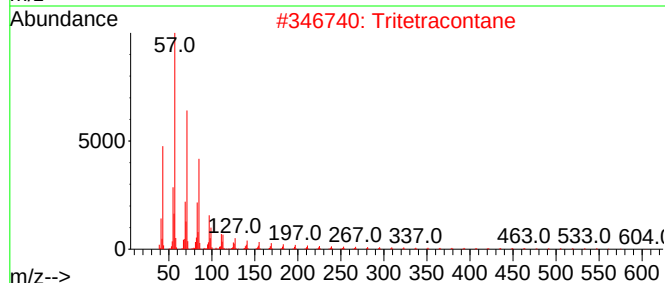
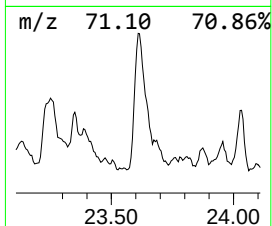
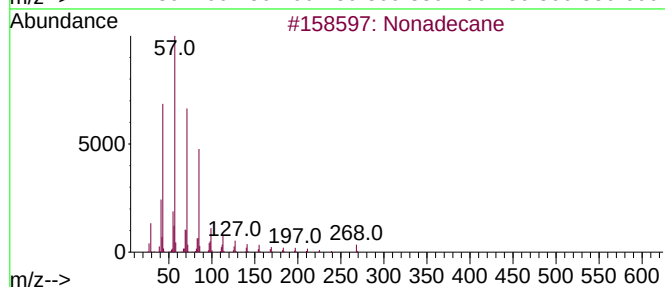
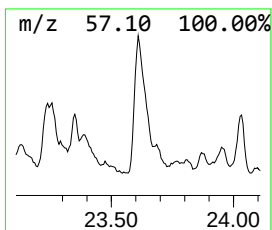
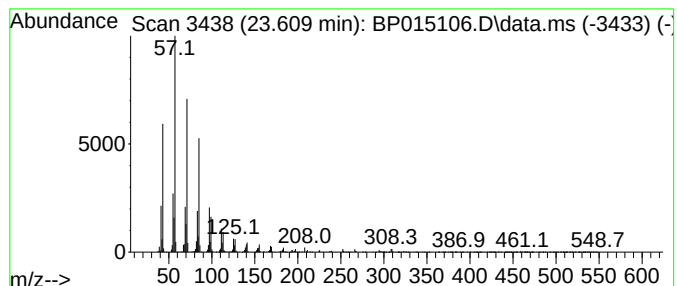
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BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.609	5.26 ng/ul	1376190	Perylene-d12	24.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	94
2	Tritetracontane	605	C43H88	007098-21-7	91
3	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
4	Hexacosane	366	C26H54	000630-01-3	83
5	Tetratetracontane	619	C44H90	007098-22-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
Data File : BP015106.D
Acq On : 16 May 2023 12:49
Operator : CG/JU
Sample : 02768-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CONN3

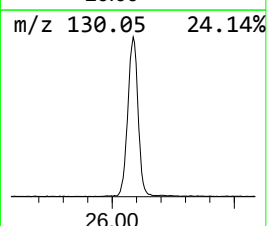
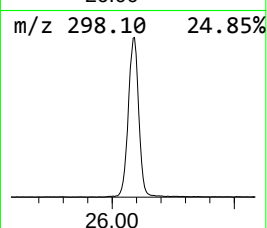
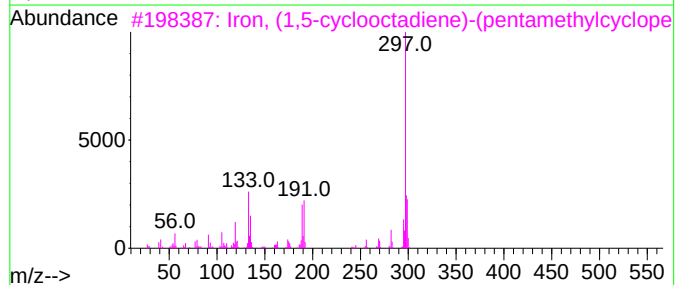
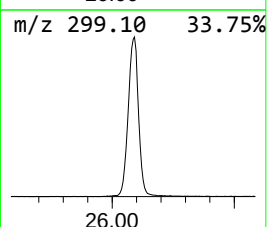
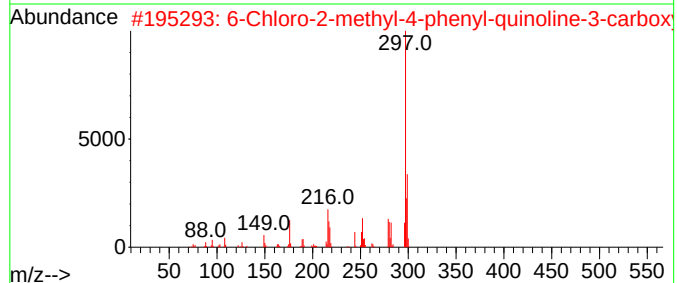
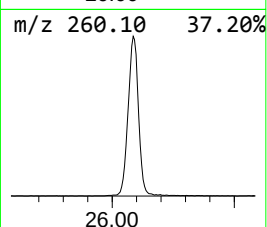
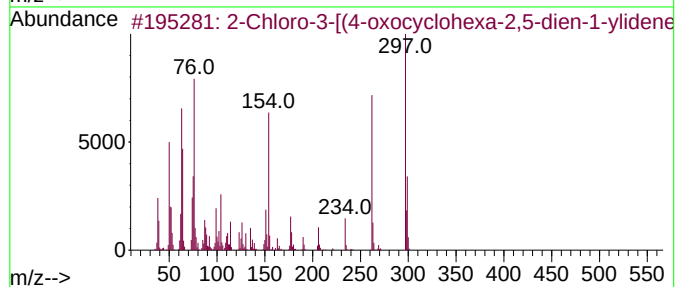
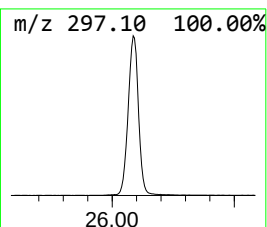
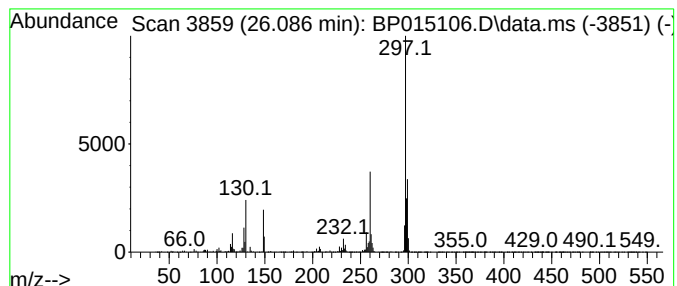
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 2-Chloro-3-[(4-oxocyclohexa... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.086	15.79 ng/ul	4129720	Perylene-d12	24.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-3-[(4-oxocyclohexa-2,5-...]	297	C16H8ClNO3	202331-78-0	53
2			6-Chloro-2-methyl-4-phenyl-quinol...	297	C17H12ClNO2	312758-85-3	52
3			Iron, (1,5-cyclooctadiene)-(pent...	299	C18H27Fe	1000161-39-5	50
4			2-[p-Chlorophenyl]-8-methylcinch...	297	C17H12ClNO2	107027-43-0	49
5			Phthalazine-1,4(2H,3H)-dione, 2-...	297	C15H11N3O4	1000272-77-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
 Data File : BP015106.D
 Acq On : 16 May 2023 12:49
 Operator : CG/JU
 Sample : 02768-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONN3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Butene, 2-chl...	3.628	5.2	ng/ul	516656	1	8.134	1985830	20.0
1-Butene, 2-(ch...	3.746	12.2	ng/ul	1207520	1	8.134	1985830	20.0
2-Butene, 1-chl...	3.975	5.0	ng/ul	492371	1	8.134	1985830	20.0
3-Methyl-3-chlo...	4.181	3.4	ng/ul	339704	1	8.134	1985830	20.0
1,1-Dimethyl-3-...	4.752	134.8	ng/ul	13379400	1	8.134	1985830	20.0
unknown-01	5.005	7.4	ng/ul	735313	1	8.134	1985830	20.0
Butane, 2,3-dic...	5.052	186.7	ng/ul	18541000	1	8.134	1985830	20.0
unknown-02	5.263	10.3	ng/ul	1026190	1	8.134	1985830	20.0
unknown-03	6.357	19.4	ng/ul	1929360	1	8.134	1985830	20.0
unknown-04	7.534	25.9	ng/ul	2573660	1	8.134	1985830	20.0
unknown-05	7.840	15.5	ng/ul	1540500	1	8.134	1985830	20.0
unknown-06	8.193	13.3	ng/ul	1315170	1	8.134	1985830	20.0
Butanoic acid, ...	8.504	5.9	ng/ul	589805	1	8.134	1985830	20.0
Propane, 1,3-di...	8.634	4.8	ng/ul	479056	1	8.134	1985830	20.0
2,2-Bis(chlorom...	8.740	3.4	ng/ul	338425	1	8.134	1985830	20.0
unknown-07	9.416	17.1	ng/ul	1696810	1	8.134	1985830	20.0
Hexanoic acid, ...	9.504	45.4	ng/ul	4512190	1	8.134	1985830	20.0
2(1H)-Pyridinon...	10.722	3.7	ng/ul	540815	2	10.963	2919220	20.0
unknown-08	11.116	4.2	ng/ul	611175	2	10.963	2919220	20.0
unknown-09	11.992	3.8	ng/ul	560873	2	10.963	2919220	20.0
unknown-10	12.639	6.8	ng/ul	991092	2	10.963	2919220	20.0
2,6-Dichloro-3,...	13.192	8.3	ng/ul	1710140	3	14.775	4125840	20.0
unknown-11	17.763	8.8	ng/ul	2182700	4	17.527	4964150	20.0
unknown-12	19.127	5.7	ng/ul	1423840	4	17.527	4964150	20.0
Adipic acid, oc...	20.898	3.7	ng/ul	944295	5	21.604	5053700	20.0
Benzoic acid, 2...	21.021	5.6	ng/ul	1406510	5	21.604	5053700	20.0
unknown-13	21.086	4.5	ng/ul	1125250	5	21.604	5053700	20.0
(DEL) Alkane: S...	21.886	2.3	ng/ul	573430	5	21.604	5053700	20.0
(DEL) Alkane: S...	21.951	2.7	ng/ul	672100	5	21.604	5053700	20.0
2-Chloro-4,6-bi...	22.104	4.7	ng/ul	1185930	5	21.604	5053700	20.0
(DEL) Alkane: S...	22.198	4.9	ng/ul	1238640	5	21.604	5053700	20.0
Heptacosane, 1-...	22.480	11.7	ng/ul	2955400	5	21.604	5053700	20.0
(DEL) Alkane: S...	23.609	5.3	ng/ul	1376190	6	24.174	5230600	20.0
2-Chloro-3-[(4-...	26.086	15.8	ng/ul	4129720	6	24.174	5230600	20.0