

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009985.D
 Acq On : 20 Apr 2022 20:05
 Operator : CG/JU
 Sample : N2472-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J40

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009985.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.258	361	369	387	rBV	7421491	12346682	34.55%	9.754%
2	7.099	676	682	693	rVB	75882	141084	0.39%	0.111%
3	7.269	704	711	720	rBV	347390	563457	1.58%	0.445%
4	7.475	739	746	760	rBV	267327	451380	1.26%	0.357%
5	7.834	798	807	816	rBV	1111552	1773926	4.96%	1.401%
6	7.981	819	832	849	rVB	5004847	8720751	24.40%	6.889%
7	8.310	878	888	909	rBV	19387544	35735234	100.00%	28.231%
8	8.646	938	945	956	rBV	218854	370598	1.04%	0.293%
9	9.099	1015	1022	1035	rBV	450706	771767	2.16%	0.610%
10	9.440	1074	1080	1089	rBV	104447	177439	0.50%	0.140%
11	9.551	1092	1099	1106	rVV2	57546	93665	0.26%	0.074%
12	9.675	1113	1120	1124	rBV3	36595	62666	0.18%	0.050%
13	9.716	1124	1127	1130	rVV2	28311	50178	0.14%	0.040%
14	9.757	1130	1134	1139	rVV2	54730	101077	0.28%	0.080%
15	9.828	1139	1146	1153	rVV	271894	476630	1.33%	0.377%
16	9.893	1153	1157	1168	rVB6	19258	41391	0.12%	0.033%
17	10.369	1229	1238	1247	rBV	392150	692805	1.94%	0.547%
18	10.622	1271	1281	1296	rBV	12987025	22798473	63.80%	18.011%
19	10.751	1296	1303	1310	rVV	652816	1054660	2.95%	0.833%
20	10.881	1318	1325	1336	rBV	182137	344377	0.96%	0.272%
21	11.140	1356	1369	1380	rBV	2858744	4842452	13.55%	3.826%
22	12.181	1540	1546	1552	rVB	25159	37445	0.10%	0.030%
23	12.316	1563	1569	1577	rVB4	19515	47109	0.13%	0.037%
24	12.740	1632	1641	1643	rBV	137031	210709	0.59%	0.166%
25	12.775	1643	1647	1655	rVV	471751	770853	2.16%	0.609%
26	12.998	1679	1685	1692	rVB9	16695	36428	0.10%	0.029%
27	13.381	1742	1750	1770	rBV	1835916	2962237	8.29%	2.340%
28	13.587	1780	1785	1793	rVB	77499	122451	0.34%	0.097%
29	13.981	1842	1852	1859	rBV	1140040	1603128	4.49%	1.266%
30	14.269	1893	1901	1915	rVV	1160532	1808842	5.06%	1.429%
31	14.575	1943	1953	1961	rBV	1080615	1604558	4.49%	1.268%
32	14.763	1980	1985	1993	rBV2	62841	104408	0.29%	0.082%
33	14.898	2001	2008	2016	rBV	382113	582816	1.63%	0.460%
34	15.569	2115	2122	2135	rBV	1621467	2442862	6.84%	1.930%
35	15.681	2135	2141	2154	rVV	326233	498664	1.40%	0.394%
36	15.810	2158	2163	2174	rVB7	25806	69327	0.19%	0.055%

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37	16.351	2251	2255	2264	rVB3	41597	69748	0.20%	0.055%
38	16.633	2298	2303	2311	rVB	65201	92749	0.26%	0.073%
39	17.328	2414	2421	2427	rBV	1317397	1985763	5.56%	1.569%
40	17.422	2431	2437	2447	rVV2	1945738	2825608	7.91%	2.232%
41	18.204	2566	2570	2578	rBV	152865	223537	0.63%	0.177%
42	18.651	2642	2646	2653	rBV2	37871	53052	0.15%	0.042%
43	19.298	2753	2756	2759	rVV2	27710	37149	0.10%	0.029%
44	19.327	2759	2761	2768	rVB2	28859	36695	0.10%	0.029%
45	19.651	2810	2816	2830	rBV	2583427	3619053	10.13%	2.859%
46	20.145	2896	2900	2913	rVB10	52630	93920	0.26%	0.074%
47	20.463	2949	2954	2963	rBV	183558	254318	0.71%	0.201%
48	21.110	3059	3064	3075	rBV	1173447	1482872	4.15%	1.171%
49	21.310	3095	3098	3103	rBV	76461	113057	0.32%	0.089%
50	21.404	3109	3114	3119	rBV	1651707	2150217	6.02%	1.699%
51	22.539	3304	3307	3313	rVB	84716	100241	0.28%	0.079%
52	23.110	3400	3404	3409	rVB	138374	209341	0.59%	0.165%
53	23.698	3497	3504	3510	rBV	1505361	3007082	8.41%	2.376%
54	23.757	3511	3514	3522	rVB	205264	349196	0.98%	0.276%
55	23.857	3524	3531	3542	rVB2	1008834	2130632	5.96%	1.683%
56	24.504	3636	3641	3649	rVB	183431	366935	1.03%	0.290%
57	24.598	3651	3657	3667	rVB3	174508	408491	1.14%	0.323%
58	24.804	3688	3692	3702	rVB2	205864	450783	1.26%	0.356%
59	25.380	3785	3790	3799	rVB2	155153	360293	1.01%	0.285%
60	26.680	4003	4011	4025	rVB	551743	1647526	4.61%	1.302%

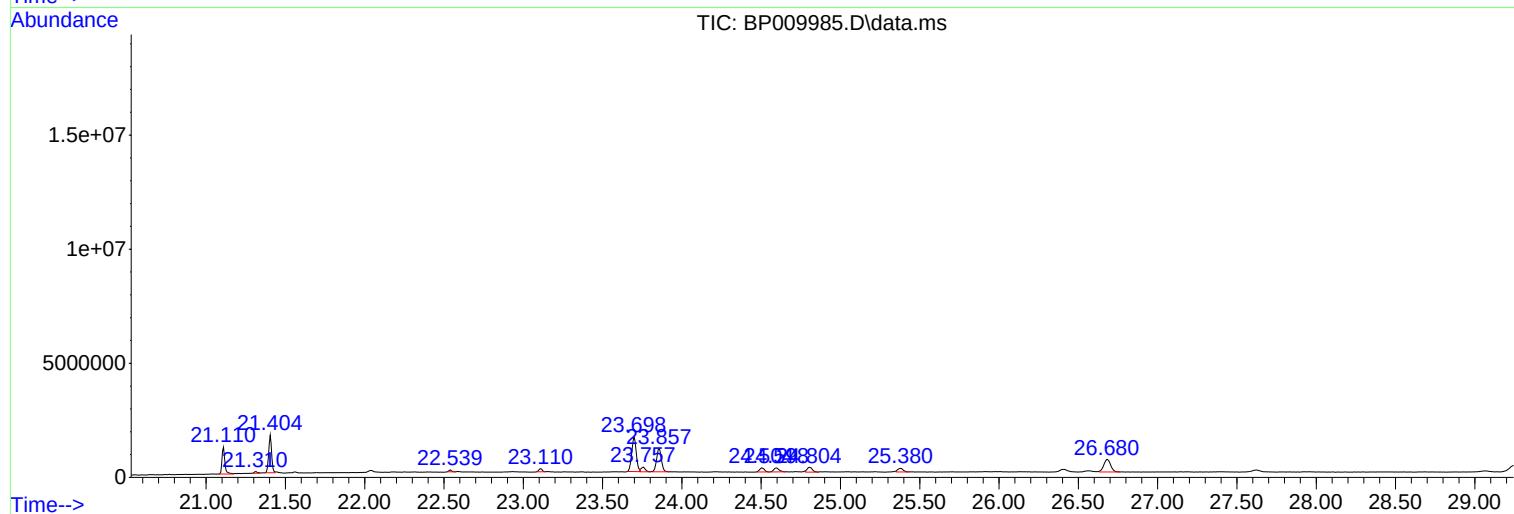
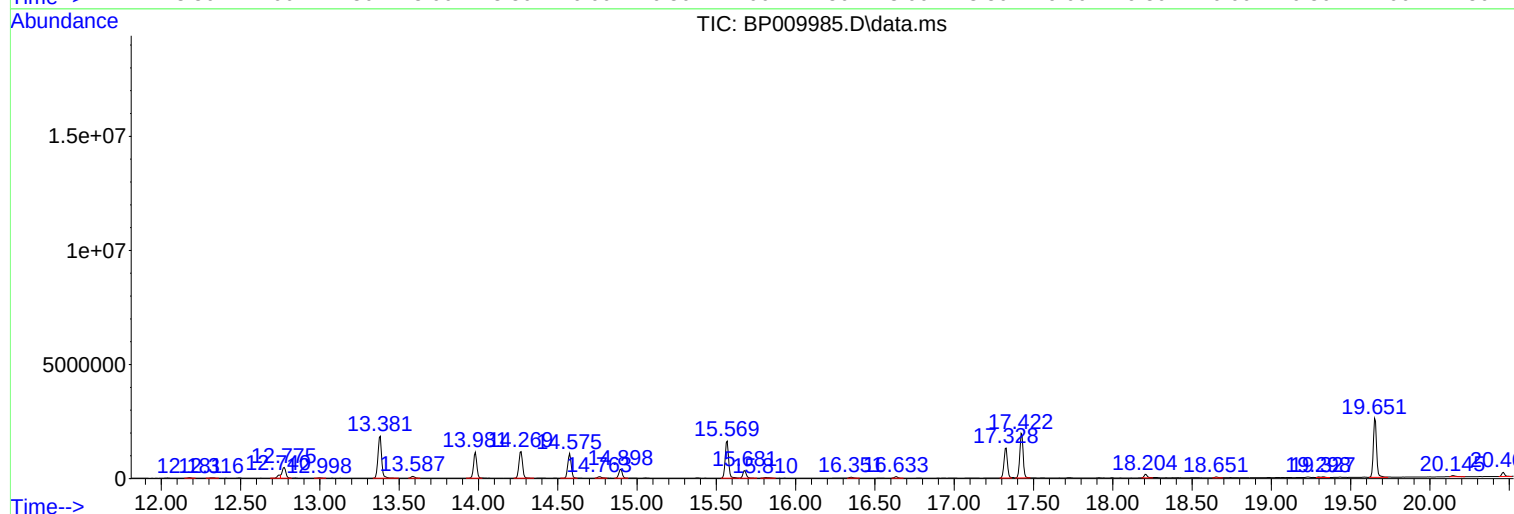
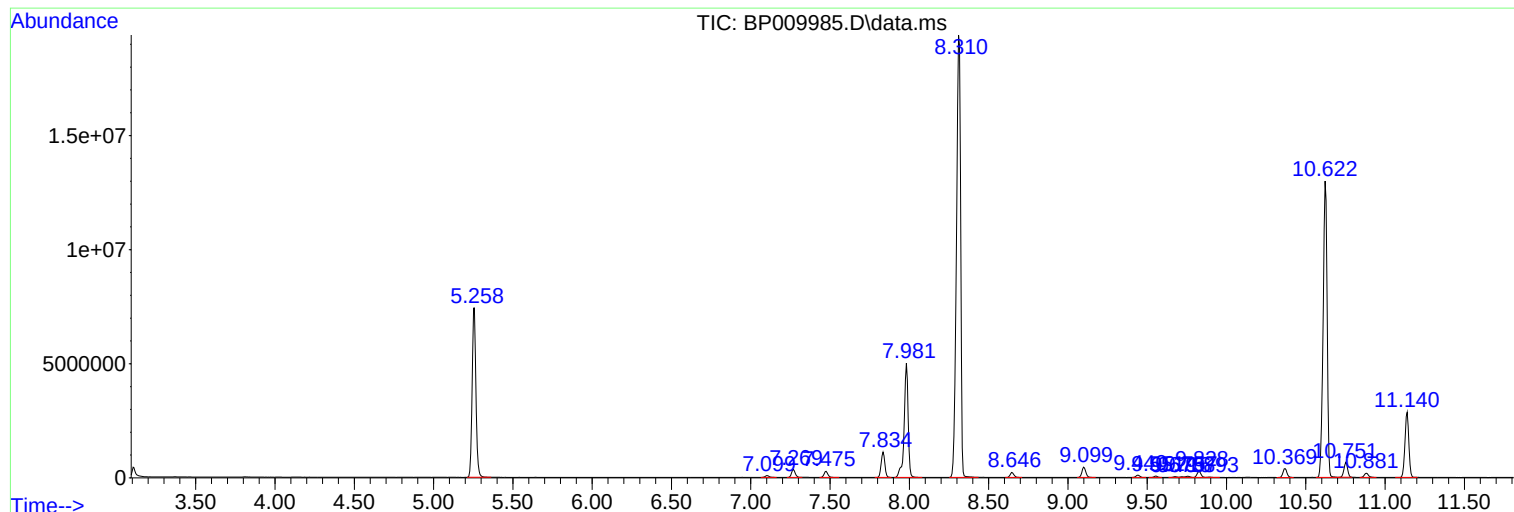
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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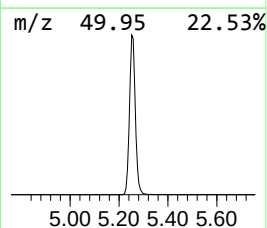
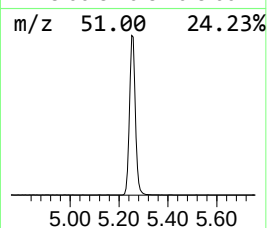
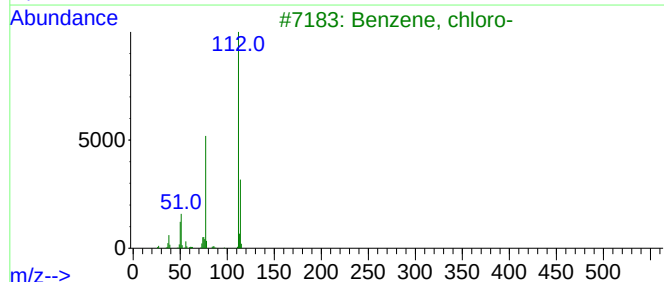
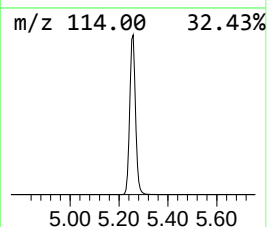
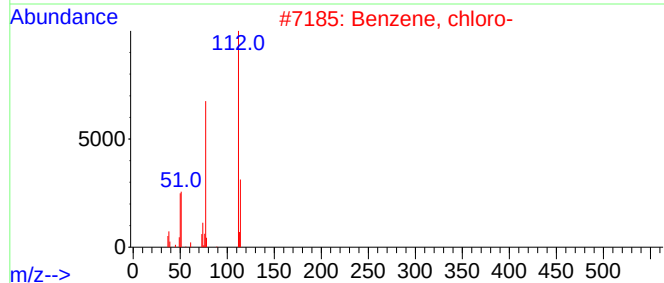
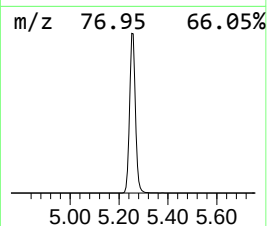
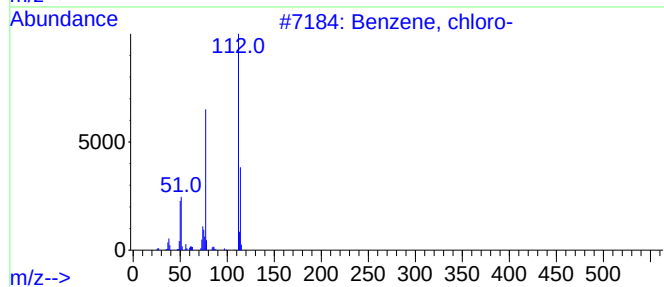
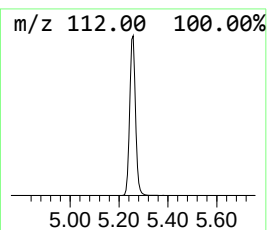
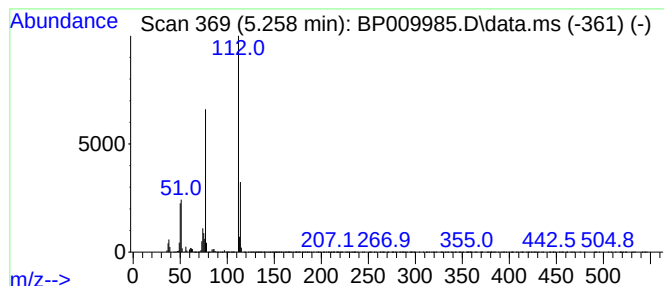
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.258	28.32 ng/ul	12346700	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



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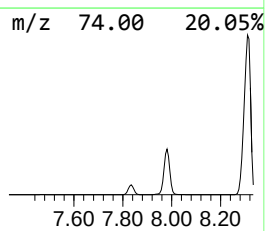
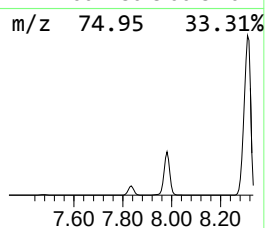
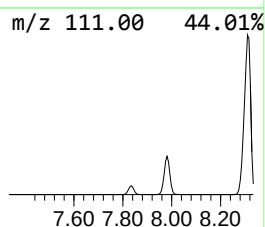
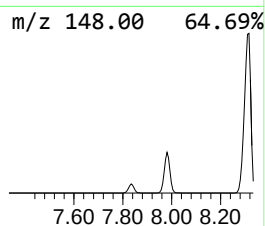
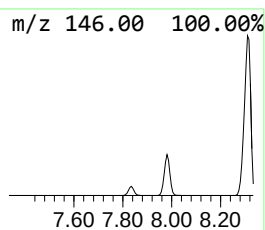
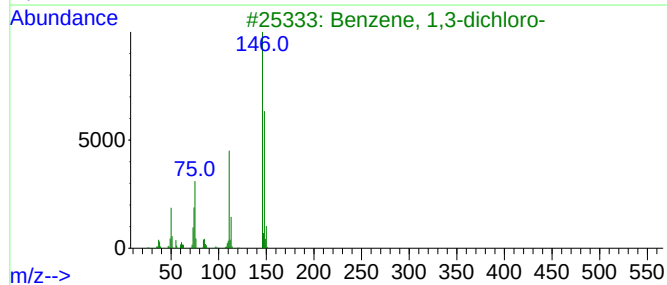
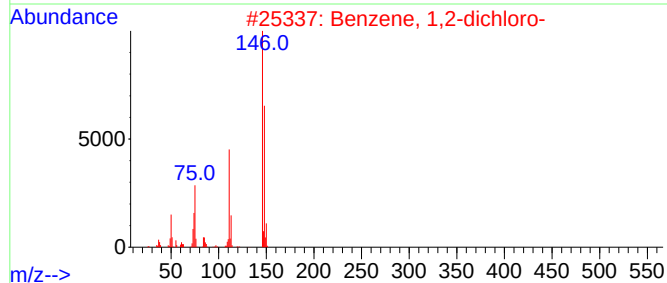
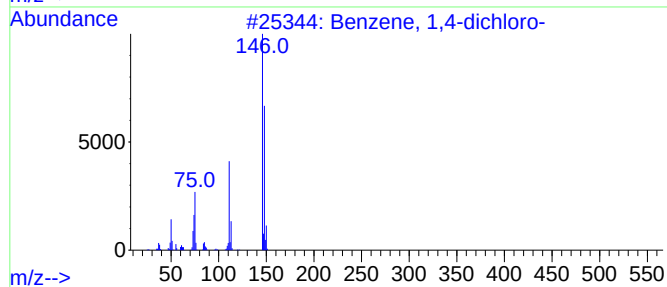
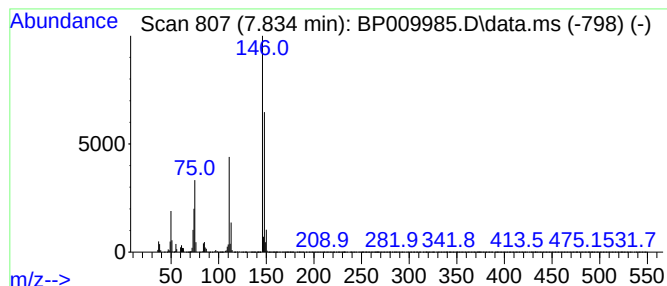
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.834	4.07 ng/ul	1773930	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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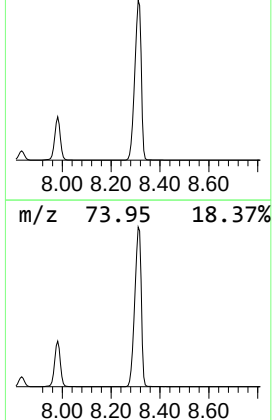
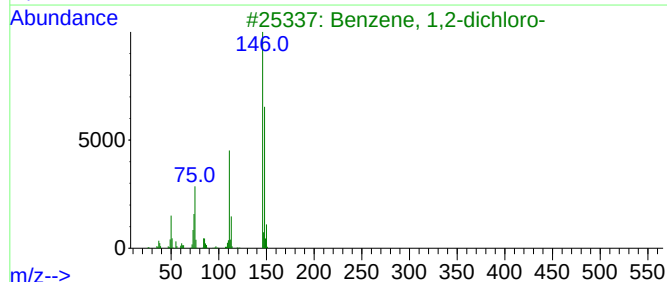
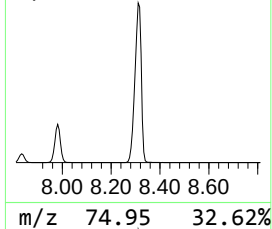
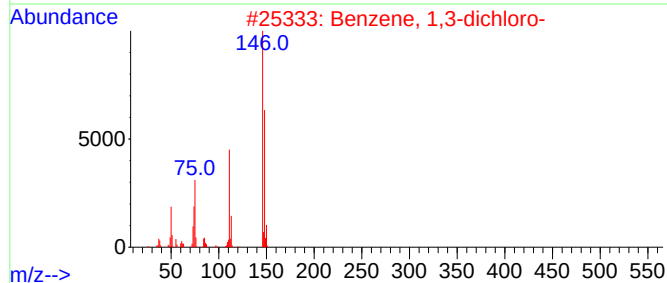
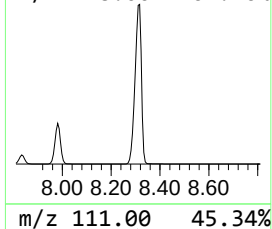
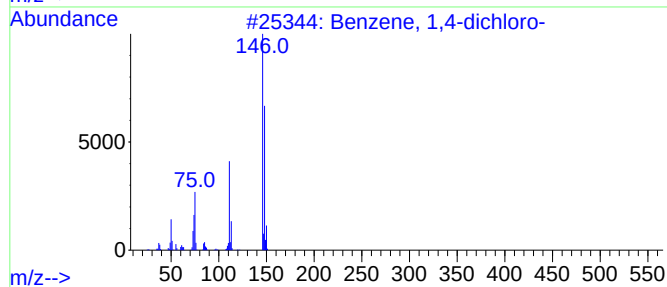
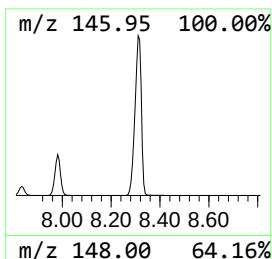
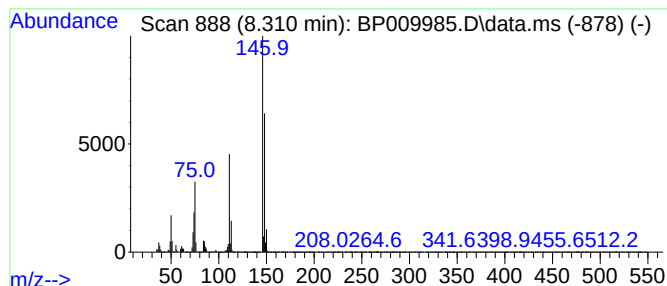
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.310	81.95 ng/ul	35735200	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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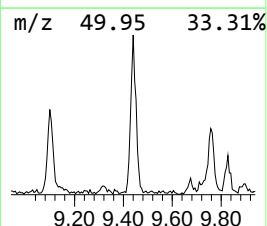
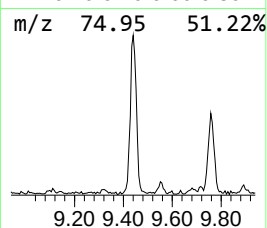
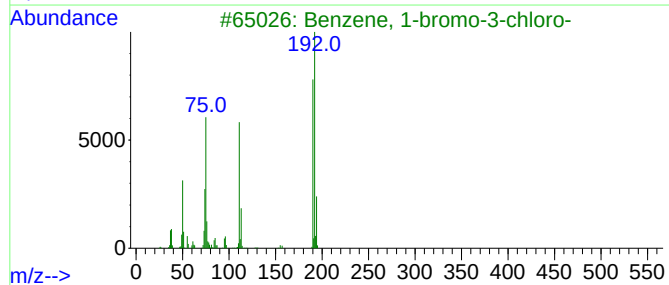
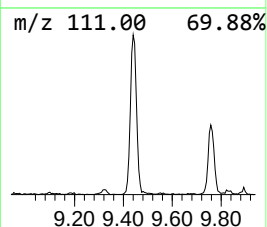
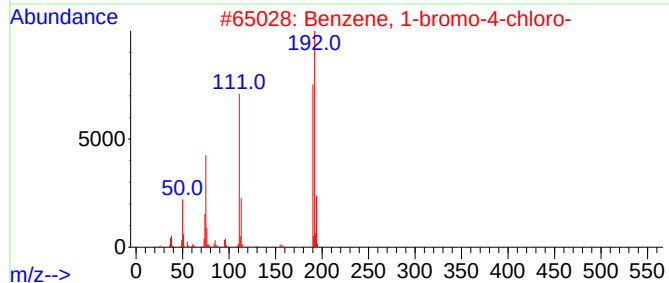
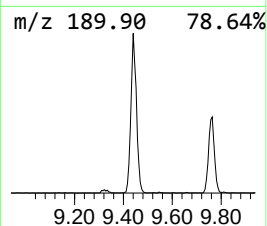
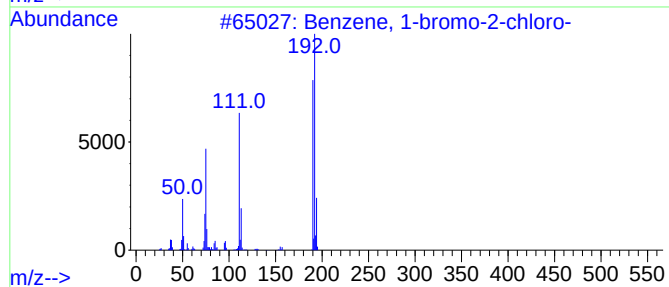
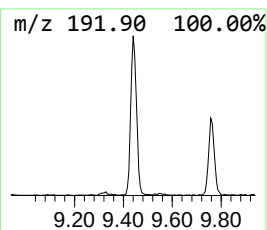
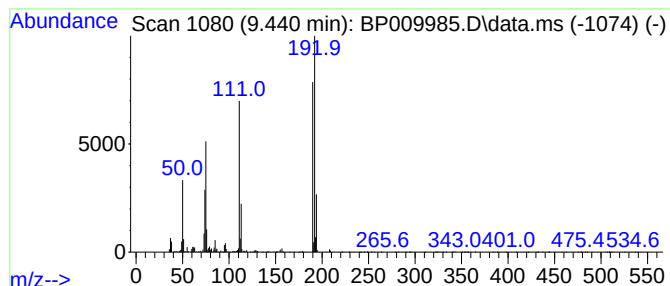
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.440	3.36 ng/ul	177439	Naphthalene-d8	10.751

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94



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C0J40

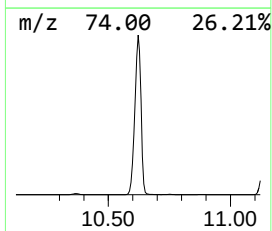
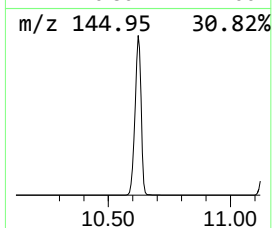
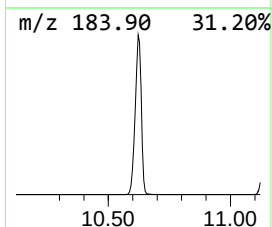
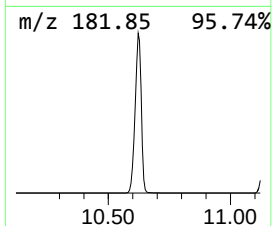
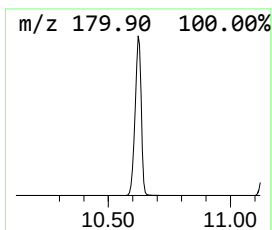
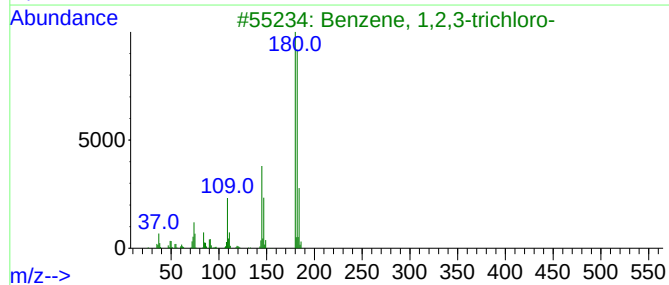
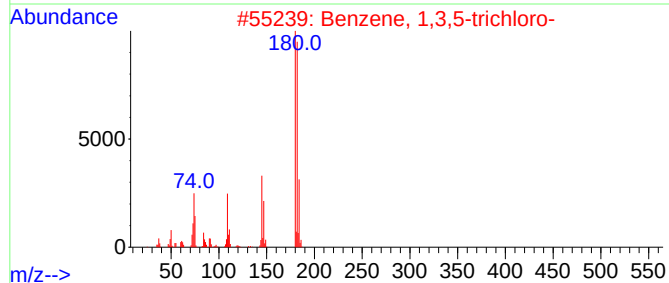
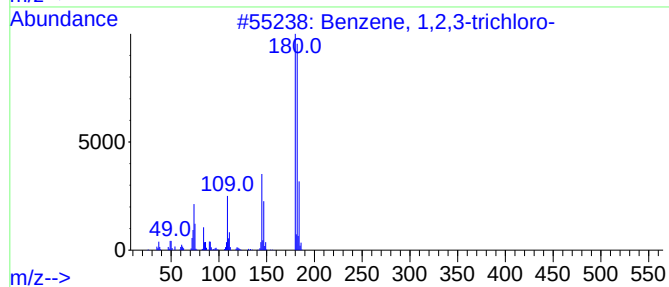
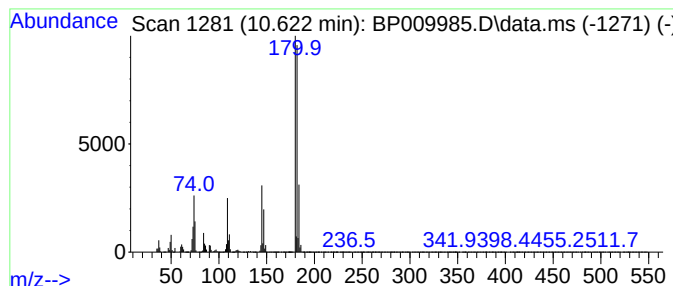
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.622	432.34 ng/ul	22798500	Naphthalene-d8	10.751

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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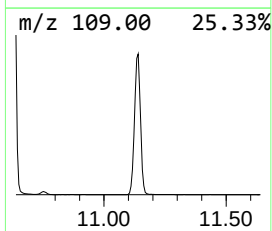
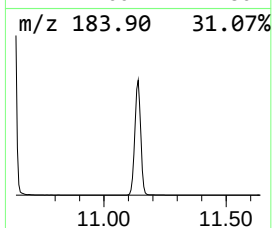
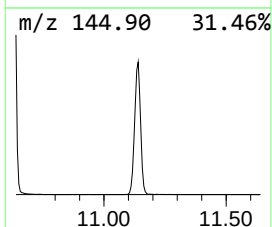
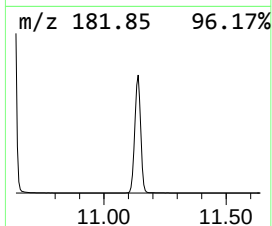
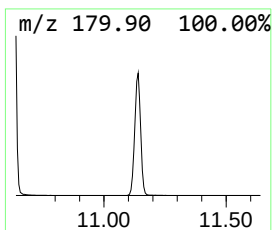
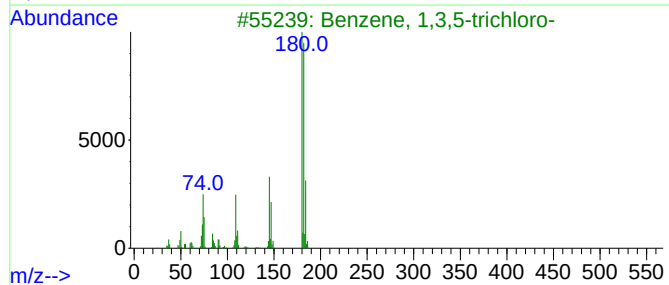
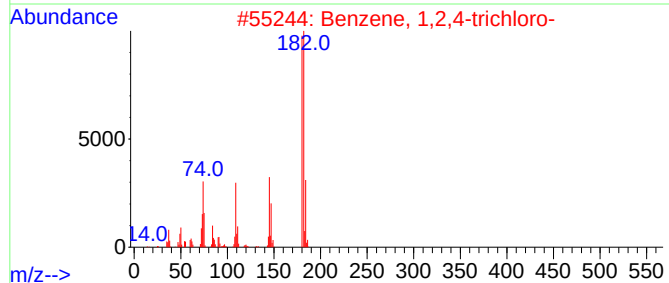
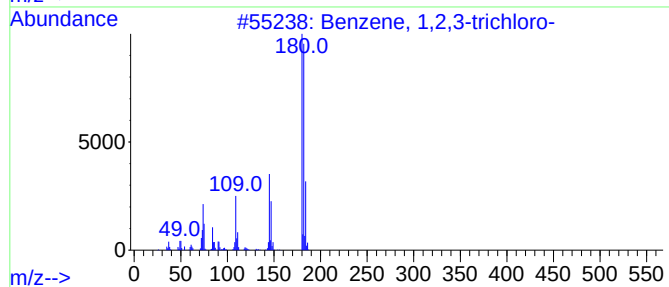
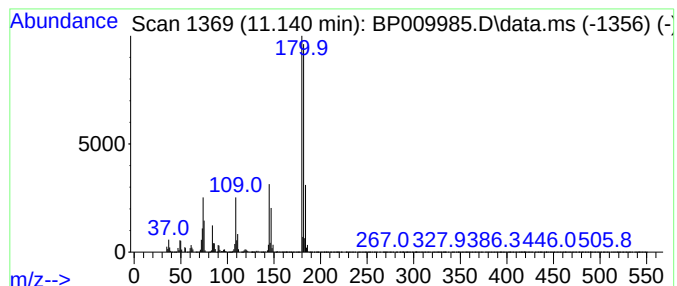
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.140	91.83 ng/ul	4842450	Naphthalene-d8	10.751

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	99
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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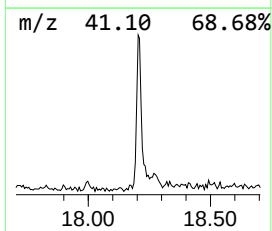
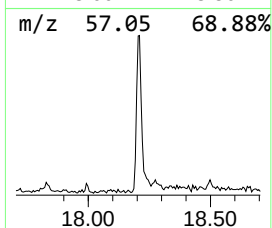
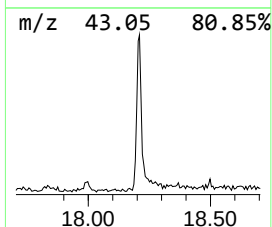
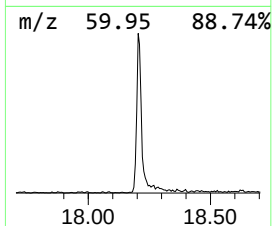
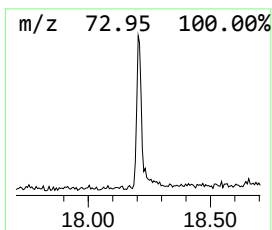
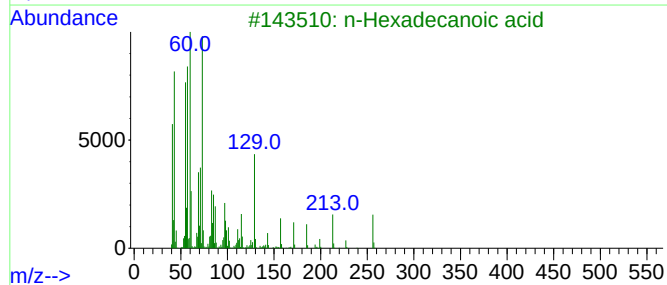
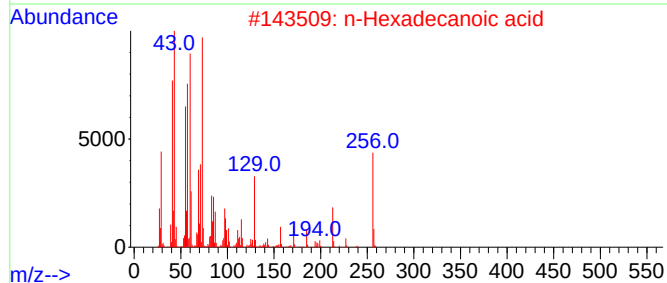
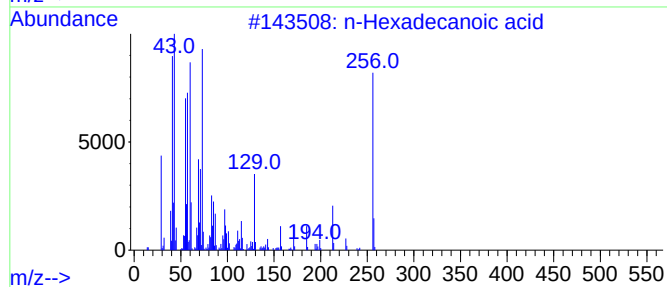
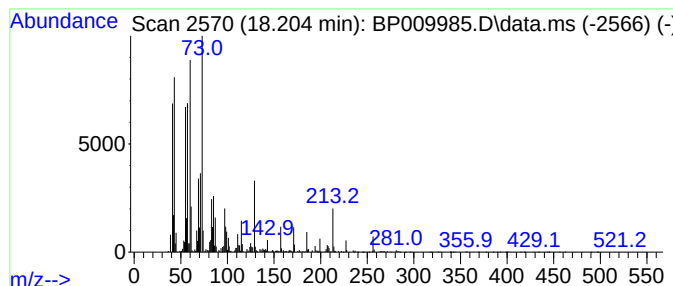
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	2.25 ng/ul	223537	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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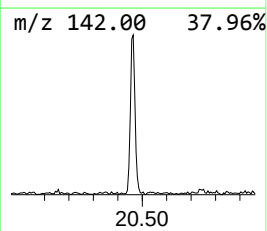
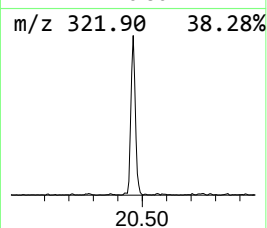
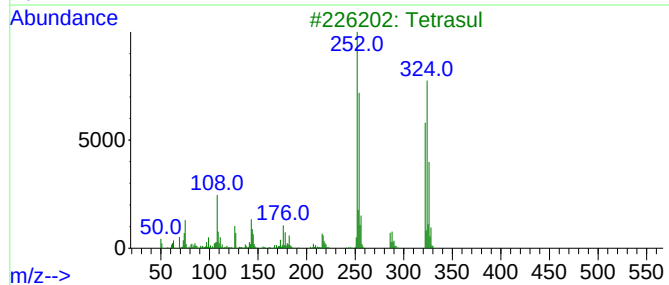
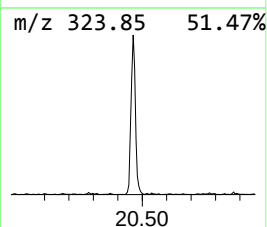
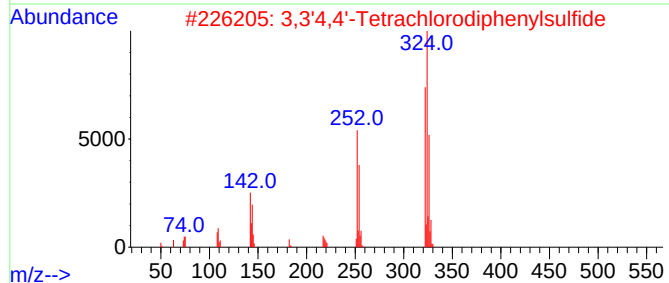
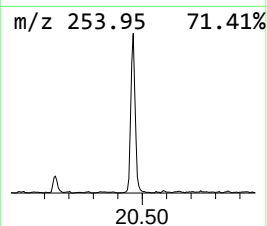
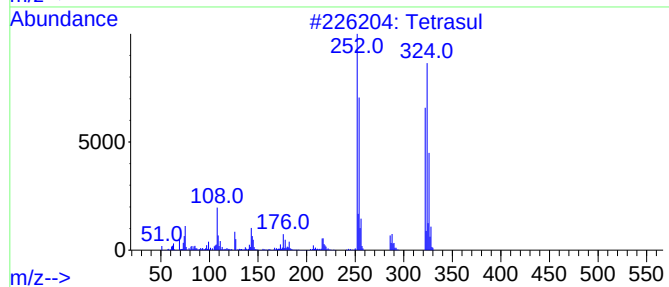
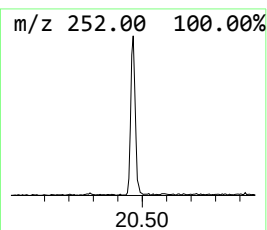
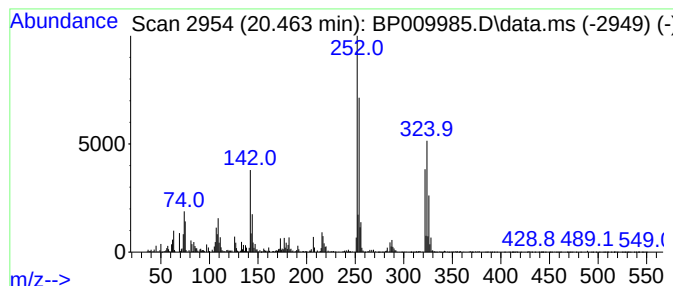
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Tetrasul Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.463	2.37 ng/ul	254318	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrasul	322	C12H6Cl4S	002227-13-6	97
2			3,3',4,4'-Tetrachlorodiphenylsulfide	322	C12H6Cl4S	076745-12-5	94
3			Tetrasul	322	C12H6Cl4S	002227-13-6	89
4			4,6'-Dihydroxy-2,2',4',6-tetrach...	322	C12H6Cl4O2	1000447-97-5	83
5			3,3'-Dichlorobenzidine, N-(2-met...	322	C16H16Cl2N2O	1000461-93-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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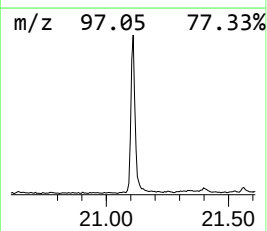
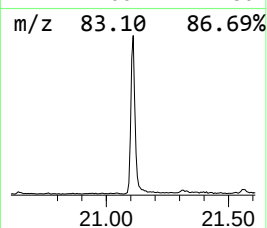
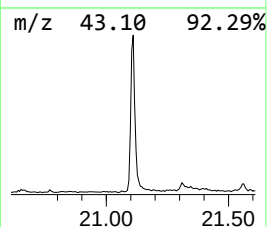
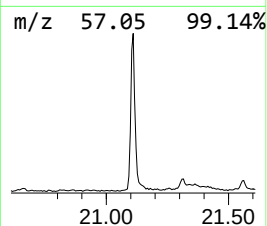
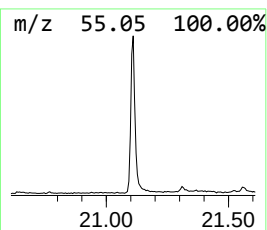
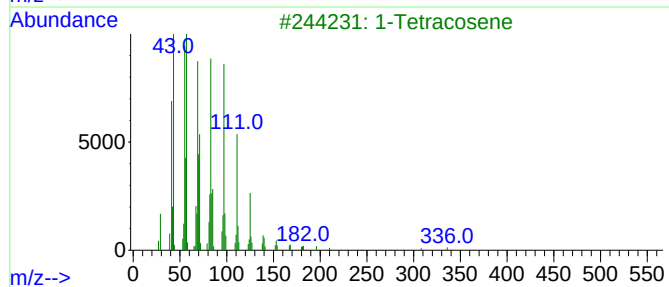
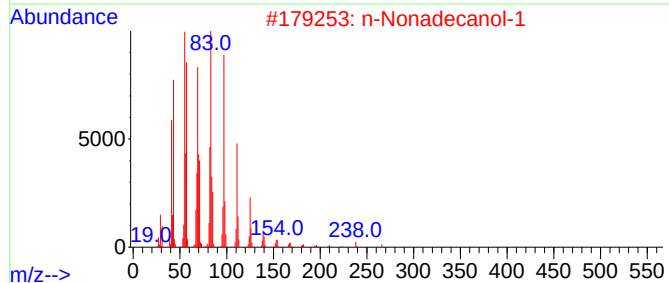
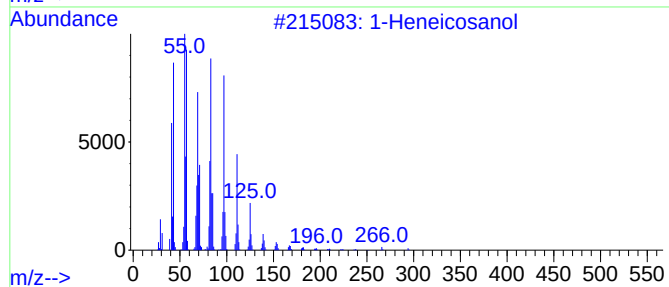
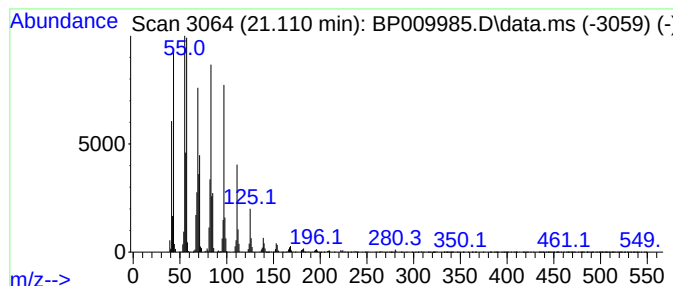
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	13.79 ng/ul	1482870	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3			1-Tetracosene	336	C24H48	010192-32-2	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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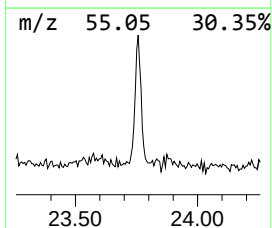
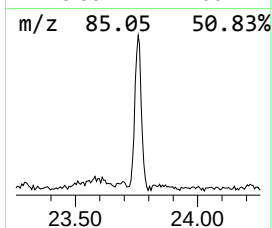
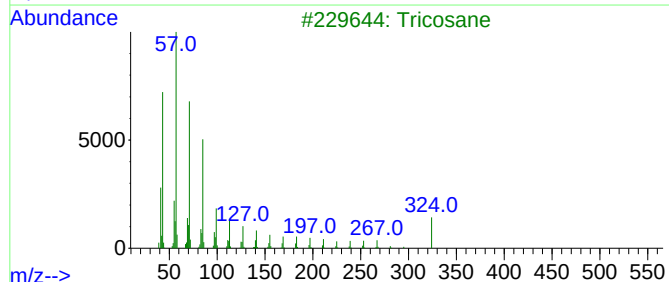
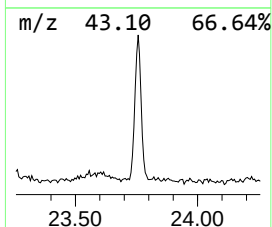
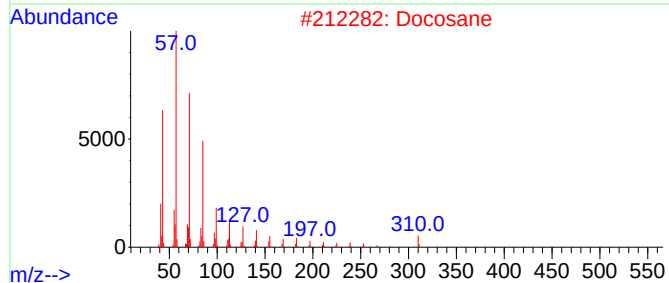
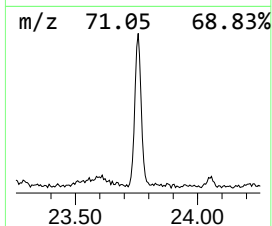
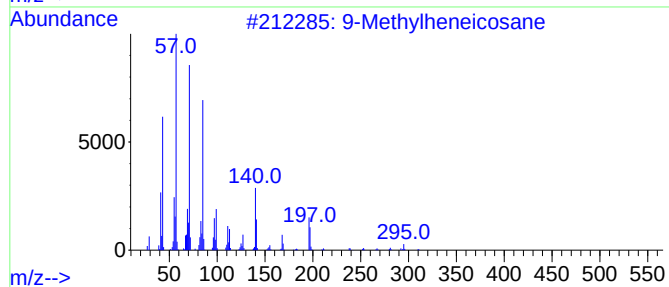
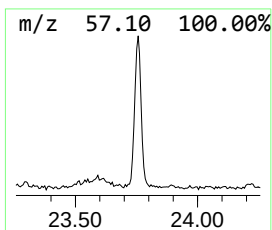
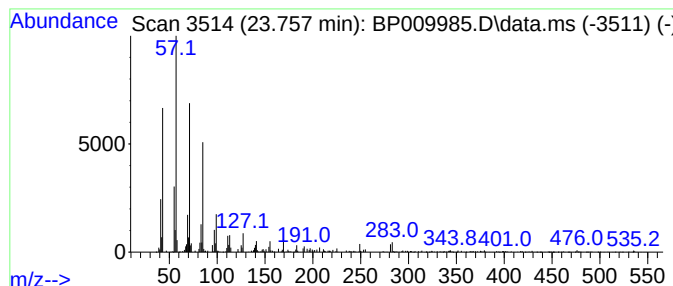
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.757	3.28 ng/ul	349196	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methylheneicosane	310	C22H46	070475-51-3	91
2			Docosane	310	C22H46	000629-97-0	91
3			Tricosane	324	C23H48	000638-67-5	90
4			2-Methylheptacosane	394	C28H58	001561-00-8	87
5			Triacontane	422	C30H62	000638-68-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
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Misc :
ALS Vial : 38 Sample Multiplier: 1

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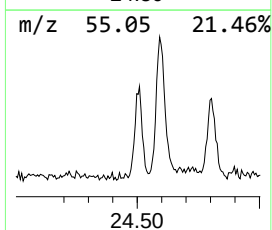
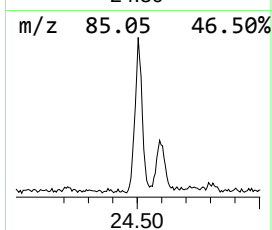
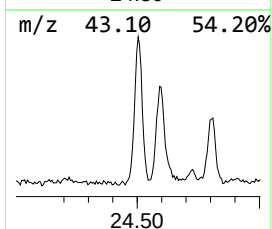
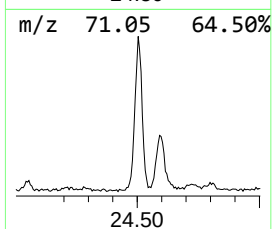
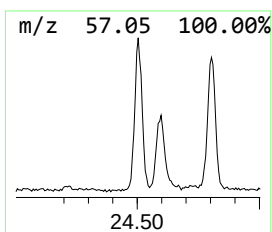
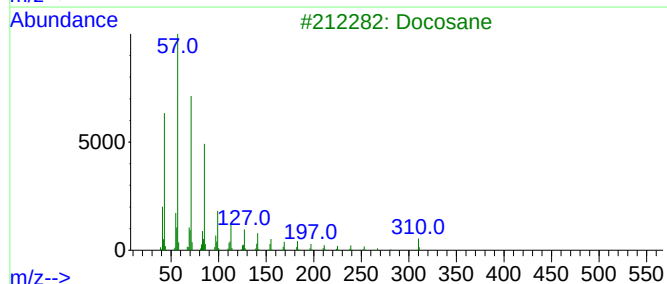
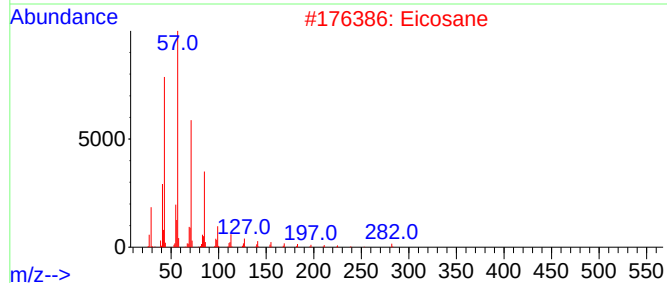
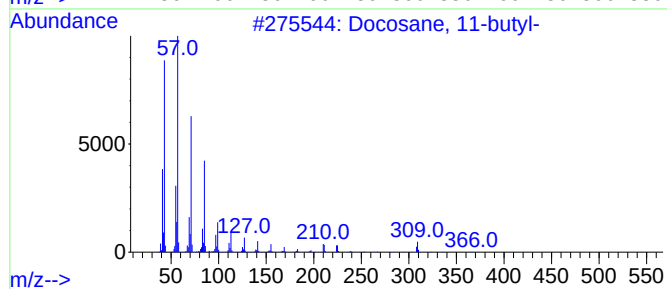
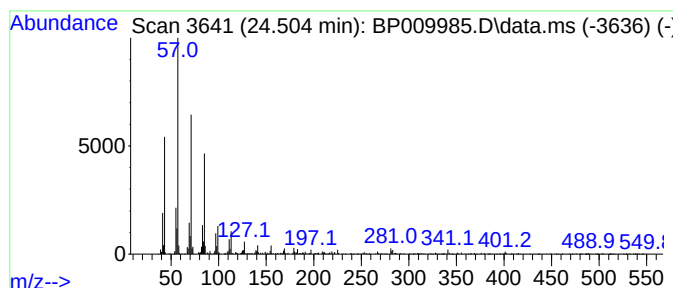
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.504	3.44 ng/ul	366935	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2			Eicosane	282	C20H42	000112-95-8	93
3			Docosane	310	C22H46	000629-97-0	93
4			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
5			Docosane, 9-butyl-	366	C26H54	055282-14-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J40

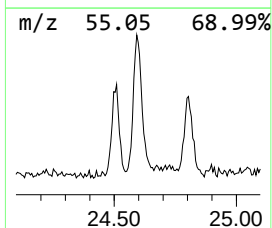
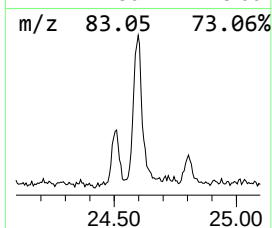
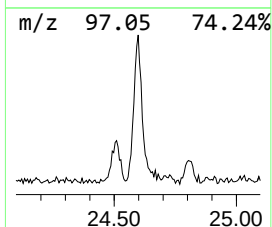
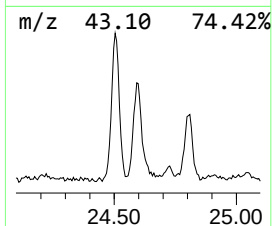
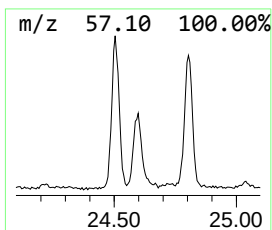
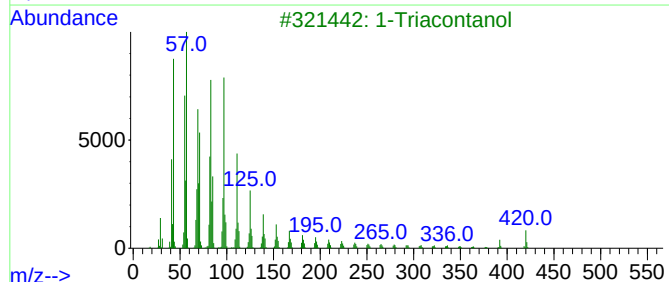
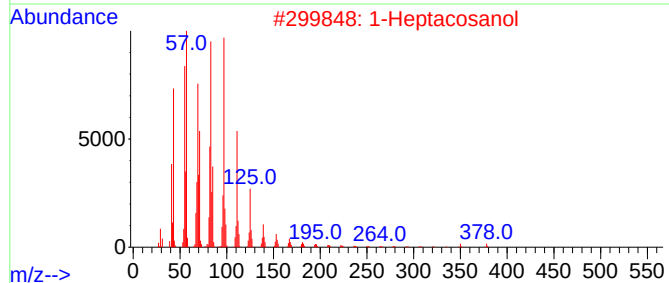
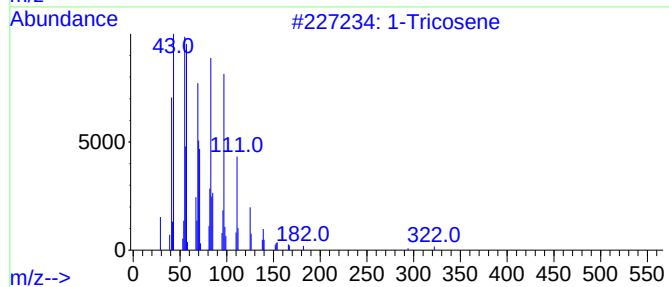
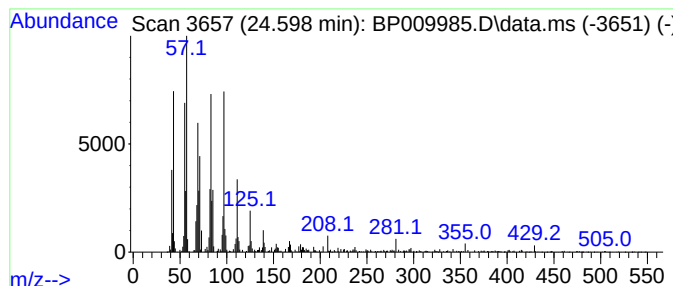
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.598	3.83 ng/ul	408491	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	95
2			1-Heptacosanol	396	C27H56O	002004-39-9	93
3			1-Triacontanol	438	C30H62O	000593-50-0	93
4			Octacosanol	410	C28H58O	000557-61-9	93
5			1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J40

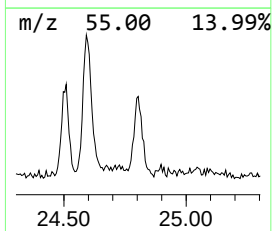
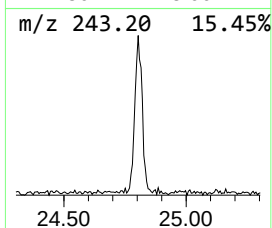
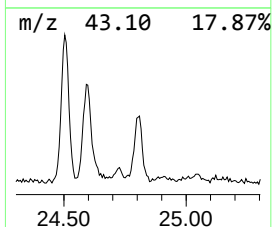
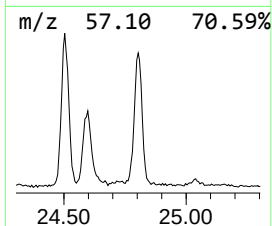
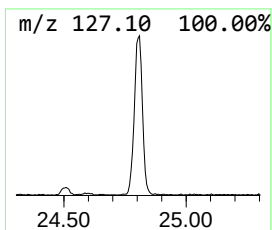
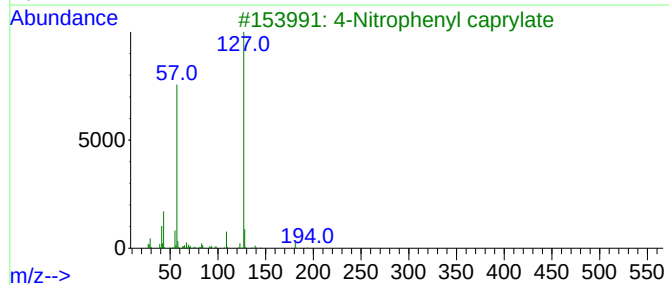
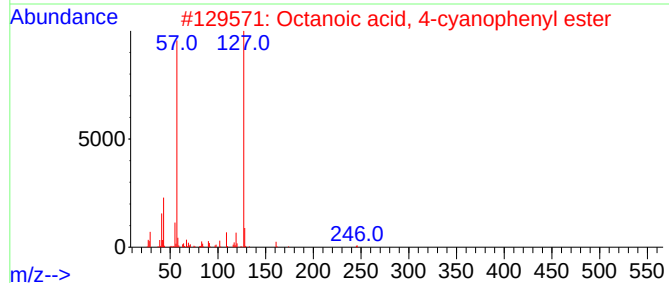
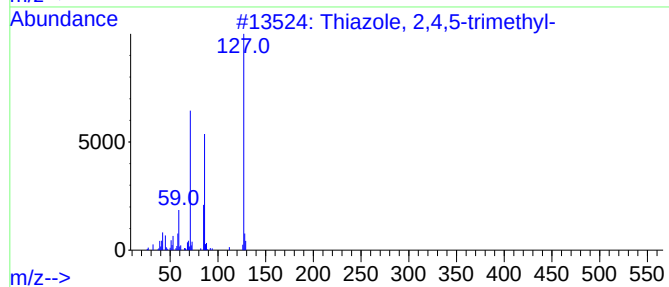
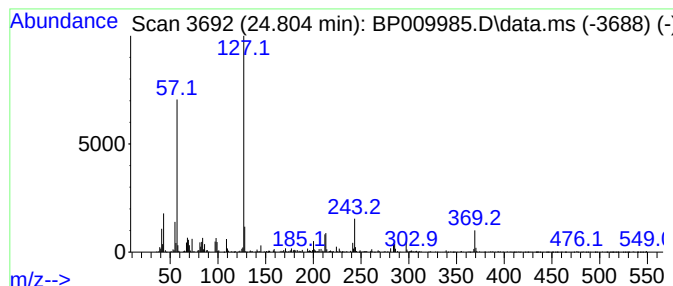
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Thiazole, 2,4,5-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.804	4.23 ng/ul	450783	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiazole, 2,4,5-trimethyl-	127	C6H9NS	013623-11-5	52
2			Octanoic acid, 4-cyanophenyl ester	245	C15H19NO2	1000307-69-8	52
3			4-Nitrophenyl caprylate	265	C14H19NO4	001956-10-1	50
4			Hexanoic acid, 2-ethyl-, anhydride	270	C16H30O3	036765-89-6	50
5			Caprylic anhydride	270	C16H30O3	000623-66-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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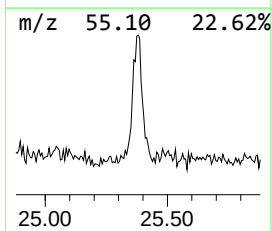
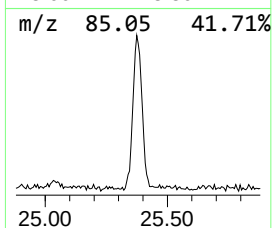
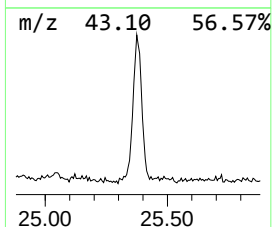
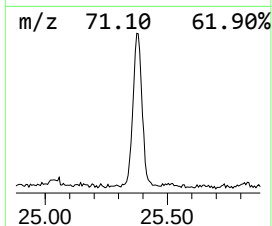
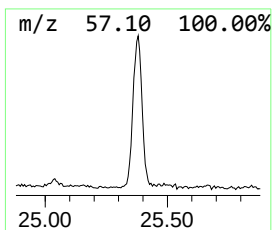
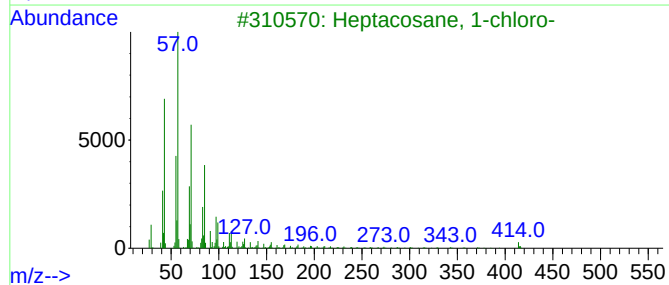
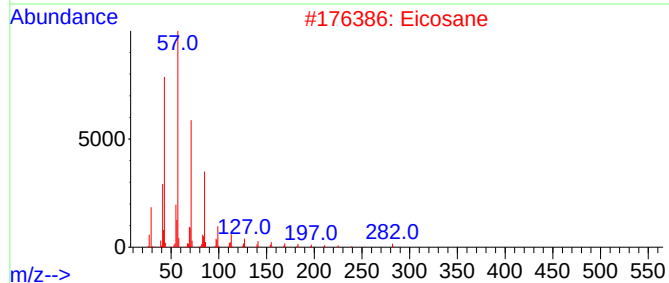
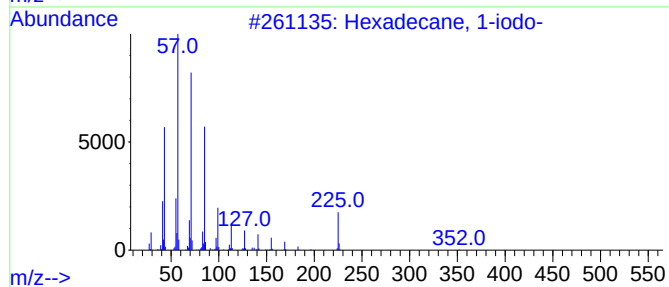
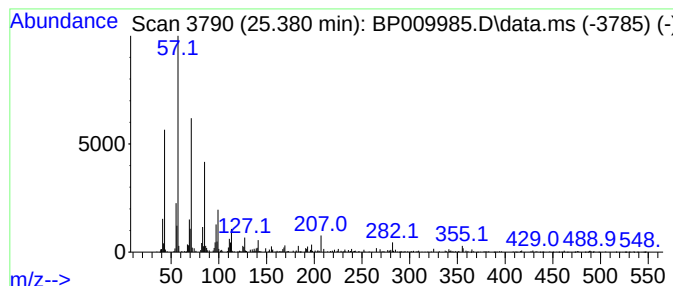
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Hexadecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.380	3.38 ng/ul	360293	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
2			Eicosane	282	C20H42	000112-95-8	91
3			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
4			Eicosane	282	C20H42	000112-95-8	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009985.D
Acq On : 20 Apr 2022 20:05
Operator : CG/JU
Sample : N2472-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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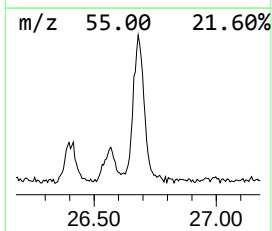
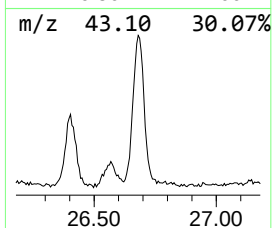
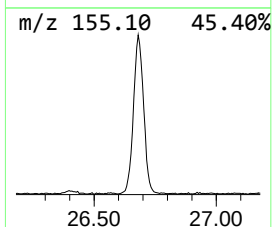
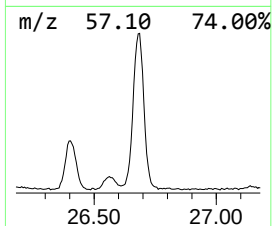
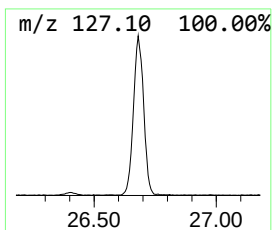
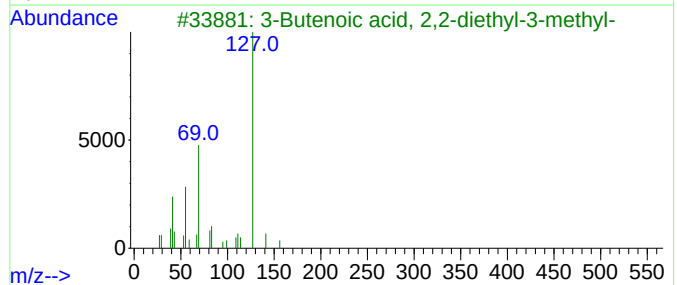
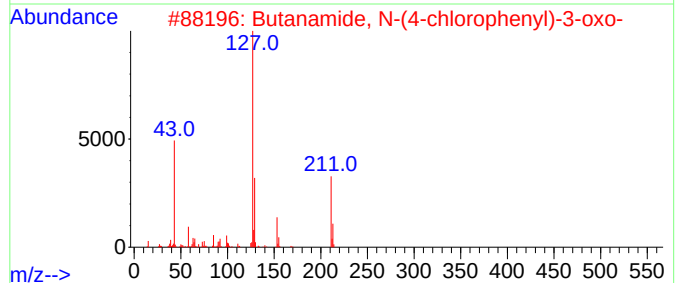
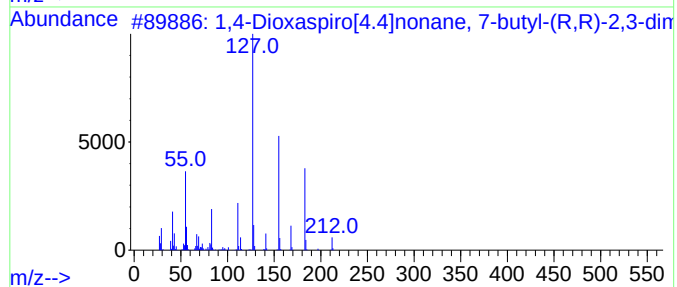
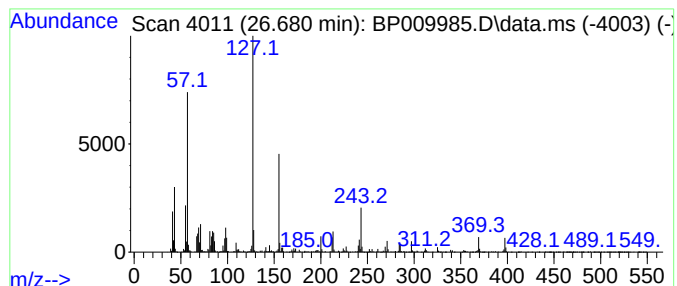
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,4-Dioxaspiro[4.4]nonane, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.680	15.47 ng/ul	1647530	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dioxaspiro[4.4]nonane, 7-but...	212	C13H24O2	104807-77-4	58
2			Butanamide, N-(4-chlorophenyl)-3...	211	C10H10ClNO2	000101-92-8	43
3			3-Butenoic acid, 2,2-diethyl-3-m...	156	C9H16O2	038477-06-4	38
4			Oxalic acid, monoamide, monohydr...	213	C10H19N3O2	1000265-20-0	38
5			2-Fluoro-5-methoxypyridine	127	C6H6FNO	136888-79-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009985.D
 Acq On : 20 Apr 2022 20:05
 Operator : CG/JU
 Sample : N2472-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
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Benzene, 1,4-di...	7.834	4.1	ng/ul	1773930	1	7.946	8720750	20.0
Benzene, 1,3-di...	8.310	82.0	ng/ul	35735200	1	7.946	8720750	20.0
Benzene, 1-brom...	9.440	3.4	ng/ul	177439	2	10.751	1054660	20.0
Benzene, 1,2,3-...	10.622	432.3	ng/ul	22798500	2	10.751	1054660	20.0
Benzene, 1,2,4-...	11.140	91.8	ng/ul	4842450	2	10.751	1054660	20.0
n-Hexadecanoic ...	18.204	2.3	ng/ul	223537	4	17.322	1985760	20.0
Tetrasul	20.463	2.4	ng/ul	254318	5	21.404	2150220	20.0
1-Heneicosanol	21.110	13.8	ng/ul	1482870	5	21.404	2150220	20.0
(DEL) Alkane: S...	23.757	3.3	ng/ul	349196	6	23.857	2130630	20.0
(DEL) Alkane: S...	24.504	3.4	ng/ul	366935	6	23.857	2130630	20.0
(DEL) Alkane: S...	24.598	3.8	ng/ul	408491	6	23.857	2130630	20.0
Thiazole, 2,4,5...	24.804	4.2	ng/ul	450783	6	23.857	2130630	20.0
Hexadecane, 1-i...	25.380	3.4	ng/ul	360293	6	23.857	2130630	20.0
1,4-Dioxaspiro[...	26.680	15.5	ng/ul	1647530	6	23.857	2130630	20.0