

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
 Data File : BP019411.D
 Acq On : 22 Jan 2024 13:34
 Operator : MA/JU
 Sample : P1142-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AZ2

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019411.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.170	11	14	18	rBV5	4455	8979	0.25%	0.023%
2	3.217	18	22	24	rVV	24628	31029	0.88%	0.079%
3	3.246	24	27	40	rVB	102744	146352	4.15%	0.372%
4	3.493	64	69	73	rBV2	7566	10942	0.31%	0.028%
5	3.640	90	94	105	rBV	136073	213532	6.06%	0.542%
6	3.799	115	121	126	rVB3	15938	25281	0.72%	0.064%
7	3.852	126	130	135	rVB3	8320	14010	0.40%	0.036%
8	3.952	143	147	151	rVB	10605	13785	0.39%	0.035%
9	4.393	218	222	229	rVB	52552	73577	2.09%	0.187%
10	4.775	279	287	291	rBV3	9532	14789	0.42%	0.038%
11	5.081	333	339	347	rVB	358862	520746	14.77%	1.323%
12	5.505	401	411	416	rVB4	13399	26312	0.75%	0.067%
13	5.711	441	446	452	rVB4	16716	27343	0.78%	0.069%
14	5.858	467	471	478	rVB7	6592	11339	0.32%	0.029%
15	6.016	493	498	502	rBV2	5430	8566	0.24%	0.022%
16	6.140	513	519	524	rBV10	4230	9141	0.26%	0.023%
17	6.263	529	540	545	rBV8	7928	21814	0.62%	0.055%
18	6.440	568	570	576	rVB5	6899	10723	0.30%	0.027%
19	6.658	600	607	614	rBV5	4511	11420	0.32%	0.029%
20	6.746	617	622	628	rVB10	4728	10717	0.30%	0.027%
21	6.963	652	659	670	rVB	120650	215235	6.11%	0.547%
22	7.110	678	684	693	rBV	488670	761317	21.60%	1.934%
23	7.205	693	700	705	rVB2	9540	18722	0.53%	0.048%
24	7.305	710	717	725	rBV	471963	744190	21.11%	1.891%
25	7.452	738	742	748	rVB3	10624	17959	0.51%	0.046%
26	7.522	750	754	760	rVB6	10352	16078	0.46%	0.041%
27	7.658	763	777	785	rBV3	43483	105347	2.99%	0.268%
28	7.769	785	796	801	rBV	518625	830662	23.57%	2.110%
29	7.869	810	813	819	rVB4	11446	18958	0.54%	0.048%
30	8.005	830	836	839	rBV8	7202	13307	0.38%	0.034%
31	8.063	840	846	850	rVV8	16208	35888	1.02%	0.091%
32	8.122	850	856	862	rVB	60444	102167	2.90%	0.260%
33	8.499	912	920	930	rVV	377245	621929	17.64%	1.580%
34	8.587	931	935	940	rVV8	6371	11653	0.33%	0.030%
35	8.681	946	951	959	rBV5	8362	21359	0.61%	0.054%
36	8.763	959	965	974	rVB3	18245	36405	1.03%	0.092%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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37	8.852	974	980	981	rBV6	7336	12144	0.34%	0.031%
38	8.940	988	995	1006	rVB	599739	988555	28.04%	2.511%
39	9.110	1019	1024	1033	rBV9	13846	36231	1.03%	0.092%
40	9.210	1040	1041	1045	rVV3	7452	8719	0.25%	0.022%

41	9.328	1052	1061	1068	rBV8	18641	40981	1.16%	0.104%
42	9.575	1096	1103	1108	rBV8	6133	13892	0.39%	0.035%
43	9.657	1108	1117	1128	rBV	511321	848262	24.06%	2.155%
44	9.752	1128	1133	1139	rBV	16171	27939	0.79%	0.071%
45	9.810	1139	1143	1145	rVV5	5489	8865	0.25%	0.023%

46	9.840	1145	1148	1152	rVB3	6894	10526	0.30%	0.027%
47	9.946	1160	1166	1169	rBV7	7076	15680	0.44%	0.040%
48	10.063	1181	1186	1190	rBV4	14884	24915	0.71%	0.063%
49	10.116	1192	1195	1202	rVB8	7635	15717	0.45%	0.040%
50	10.199	1202	1209	1220	rBV	771992	1360499	38.60%	3.456%

51	10.428	1238	1248	1260	rBV	260125	447121	12.68%	1.136%
52	10.563	1264	1271	1277	rBV	689435	1136819	32.25%	2.888%
53	10.616	1278	1280	1285	rVV4	22908	34553	0.98%	0.088%
54	10.663	1286	1288	1291	rVV3	7220	11090	0.31%	0.028%
55	10.716	1291	1297	1310	rVV	556150	975533	27.68%	2.478%

56	10.946	1331	1336	1340	rBV2	32313	59849	1.70%	0.152%
57	11.122	1359	1366	1376	rVB	64184	121922	3.46%	0.310%
58	11.281	1388	1393	1398	rVB9	10038	18746	0.53%	0.048%
59	11.381	1406	1410	1416	rBV9	6835	15004	0.43%	0.038%
60	11.463	1422	1424	1433	rVB9	8382	16023	0.45%	0.041%

61	11.598	1442	1447	1450	rBV3	14467	23779	0.67%	0.060%
62	11.681	1457	1461	1469	rVB7	9945	18343	0.52%	0.047%
63	11.928	1492	1503	1510	rBV	90083	166911	4.74%	0.424%
64	12.040	1518	1522	1526	rBV3	8529	12756	0.36%	0.032%
65	12.122	1534	1536	1539	rBV3	8610	10322	0.29%	0.026%

66	12.157	1539	1542	1545	rVB3	9792	11247	0.32%	0.029%
67	12.240	1552	1556	1562	rVB8	8489	14416	0.41%	0.037%
68	12.522	1601	1604	1607	rVB5	10890	13258	0.38%	0.034%
69	12.669	1625	1629	1632	rBV	21166	24834	0.70%	0.063%
70	12.763	1641	1645	1648	rBV6	10396	15929	0.45%	0.040%

71	12.934	1672	1674	1684	rVB6	8194	14348	0.41%	0.036%
72	13.151	1709	1711	1717	rVB4	6876	9642	0.27%	0.024%
73	13.234	1717	1725	1727	rBV5	21283	56534	1.60%	0.144%
74	13.263	1727	1730	1735	rVV	83701	128149	3.64%	0.326%
75	13.310	1735	1738	1746	rVB3	31639	59169	1.68%	0.150%

76	13.398	1748	1753	1755	rBV6	10696	18949	0.54%	0.048%
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 Title : SVOA CALIBRATION

77	13.745	1808	1812	1814	rBV4	12676	18621	0.53%	0.047%
78	13.834	1821	1827	1833	rBV	1667730	2258976	64.09%	5.739%
79	13.887	1833	1836	1844	rVB9	23865	51222	1.45%	0.130%
80	14.069	1864	1867	1868	rBV3	11892	13577	0.39%	0.034%
81	14.110	1868	1874	1882	rVB	1718920	2581846	73.25%	6.559%
82	14.228	1892	1894	1897	rBV4	10421	12749	0.36%	0.032%
83	14.416	1920	1926	1931	rBV	1011247	1470746	41.72%	3.736%
84	14.457	1931	1933	1940	rVB2	64146	92260	2.62%	0.234%
85	14.675	1964	1970	1977	rBV4	108474	223766	6.35%	0.568%
86	14.981	2018	2022	2026	rBV	34838	68425	1.94%	0.174%
87	15.216	2058	2062	2066	rBV	58854	72316	2.05%	0.184%
88	15.410	2089	2095	2102	rBV	2218460	3208004	91.01%	8.150%
89	15.551	2114	2119	2133	rVB	723987	994396	28.21%	2.526%
90	16.686	2308	2312	2321	rVB	101164	140603	3.99%	0.357%
91	17.169	2389	2394	2402	rVB	1163071	1704311	48.35%	4.330%
92	17.269	2406	2411	2422	rVB2	2292491	3245742	92.08%	8.245%
93	18.069	2543	2547	2554	rBV	663951	880329	24.97%	2.236%
94	19.304	2754	2757	2768	rVB	243208	312850	8.88%	0.795%
95	19.522	2788	2794	2799	rBV	2669642	3524892	100.00%	8.955%
96	20.527	2962	2965	2969	rVB	79062	82095	2.33%	0.209%
97	20.986	3039	3043	3051	rBV	466862	620345	17.60%	1.576%
98	21.274	3088	3092	3098	rBV	1215632	1505285	42.70%	3.824%
99	23.474	3459	3466	3479	rVB2	1569545	3148015	89.31%	7.997%
100	23.621	3486	3491	3501	rVB	754412	1503014	42.64%	3.818%

Sum of corrected areas: 39364129

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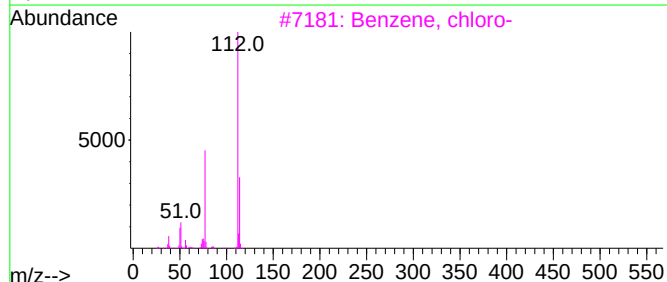
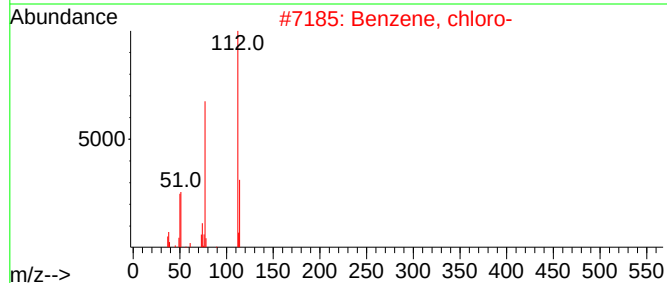
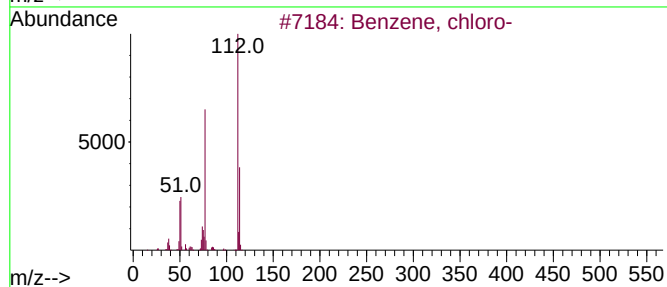
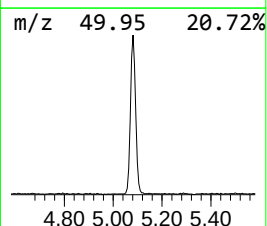
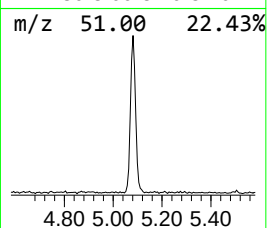
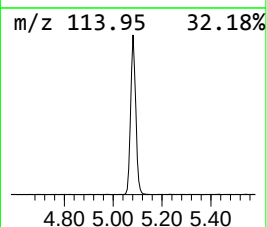
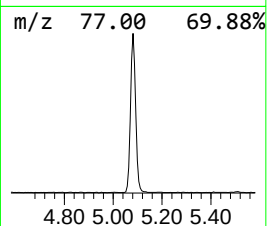
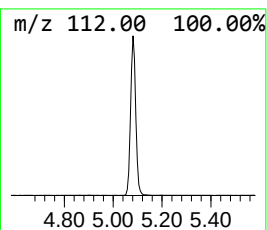
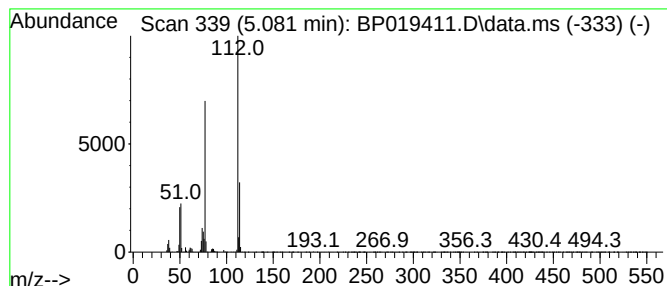
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.081	12.54 ng/ul	520746	1,4-Dichlorobenzene-d4	7.769

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	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



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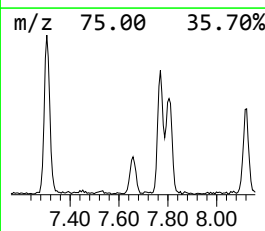
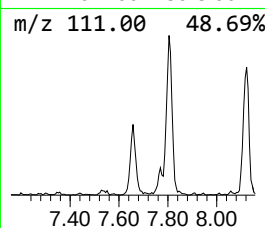
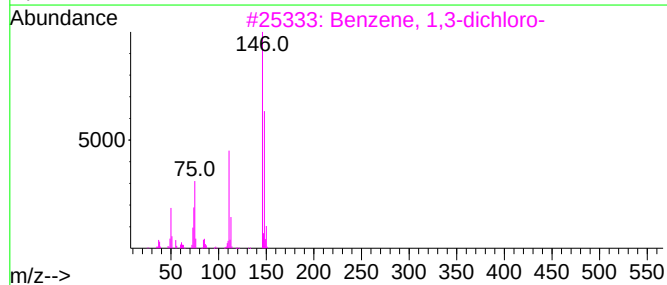
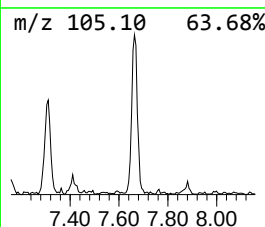
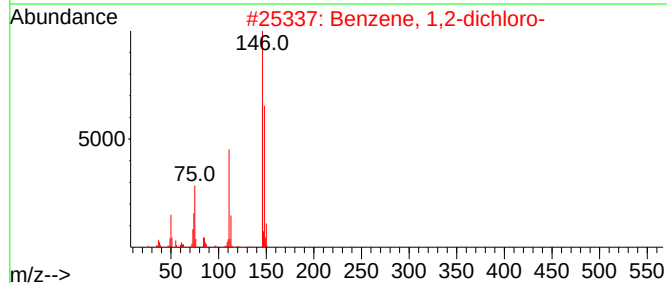
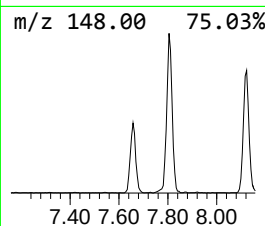
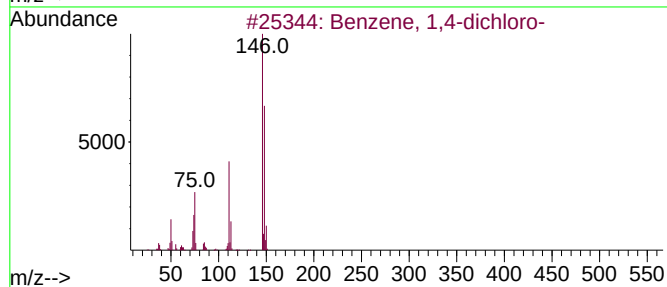
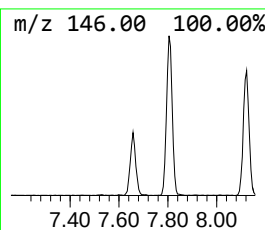
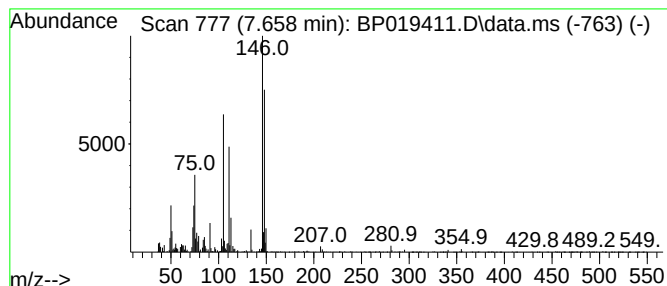
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.658	2.54 ng/ul	105347	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95



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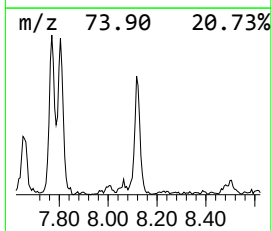
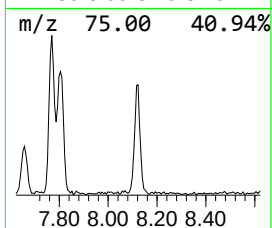
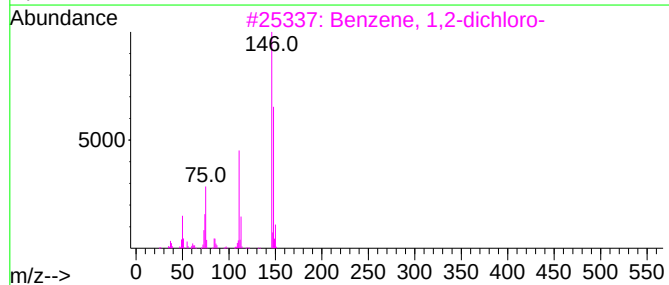
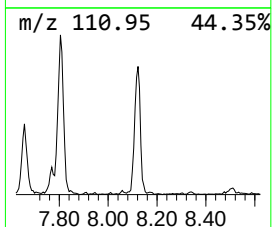
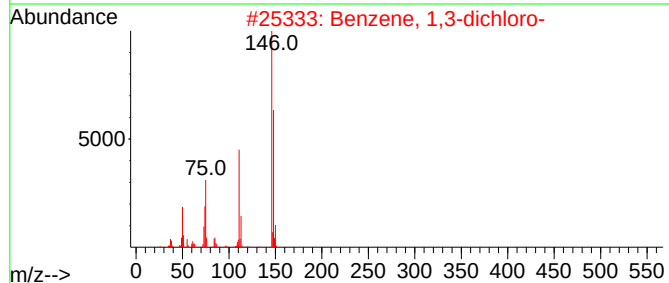
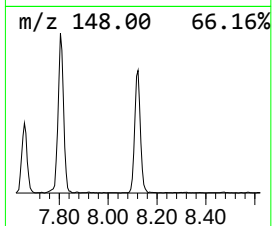
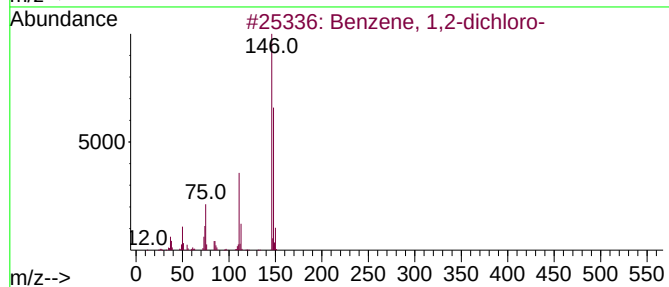
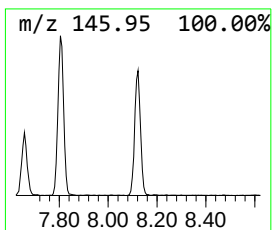
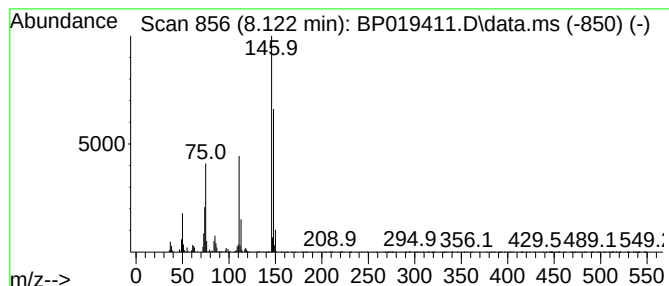
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	2.46 ng/ul	102167	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94



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ClientSampleId :
C0AZ2

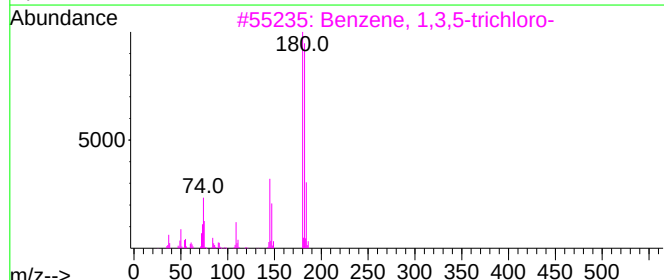
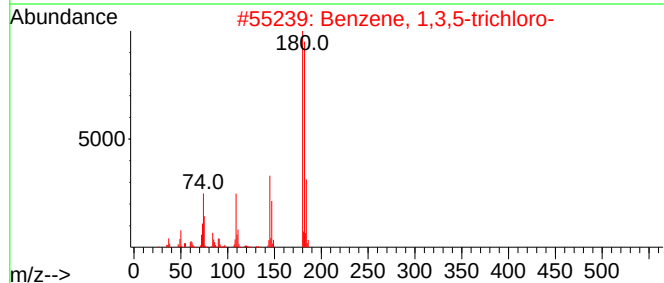
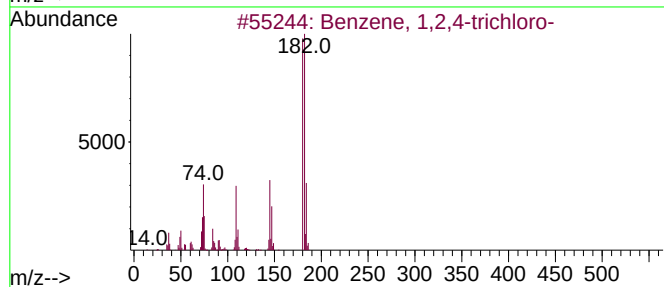
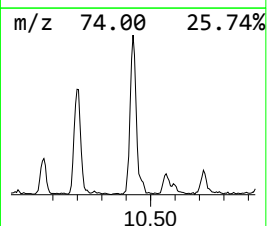
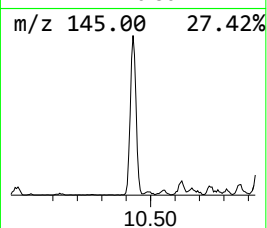
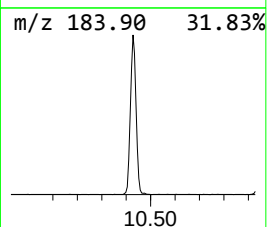
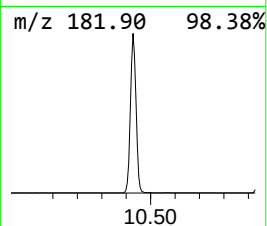
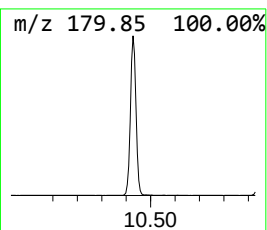
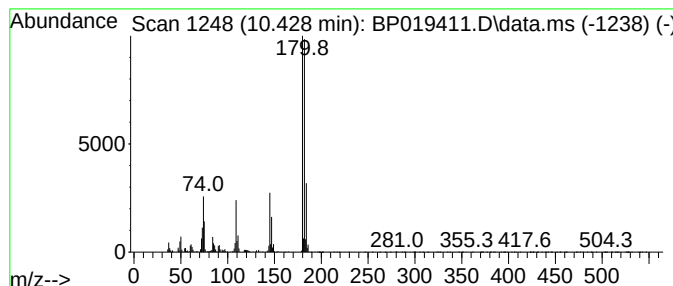
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.428	7.87 ng/ul	447121	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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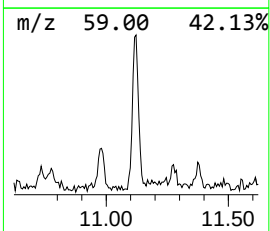
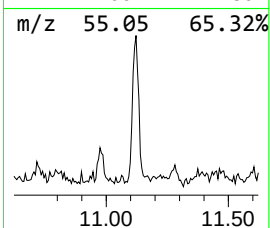
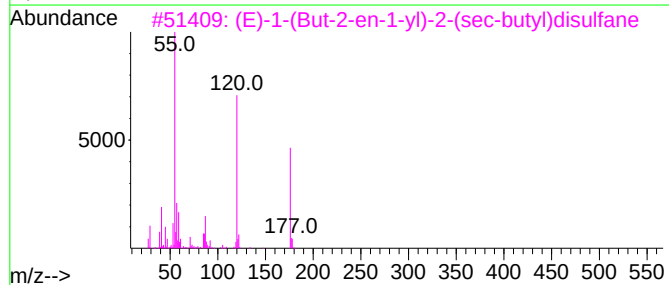
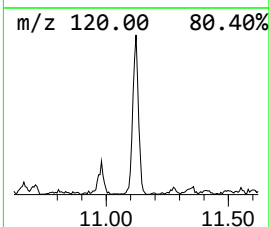
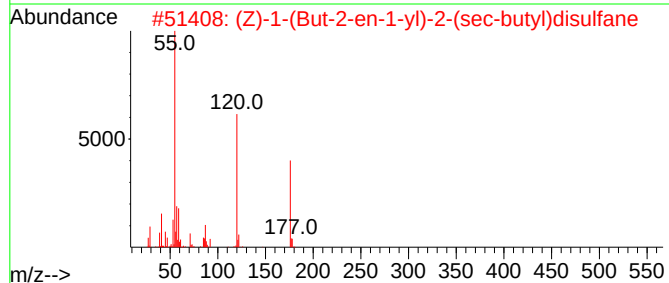
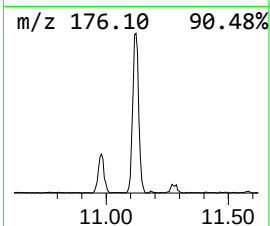
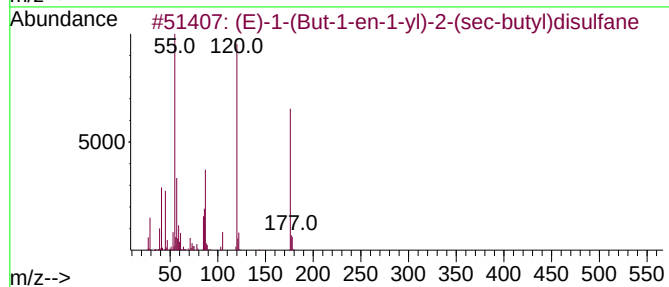
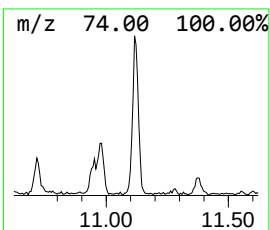
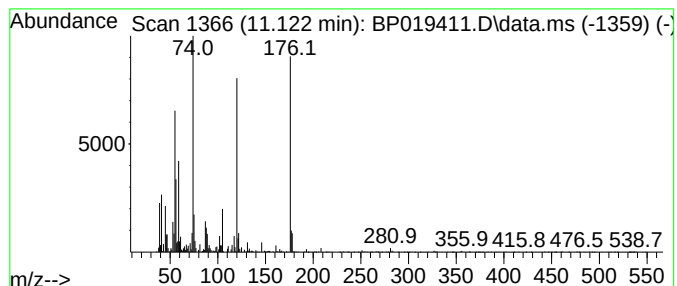
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (E)-1-(But-1-en-1-yl)-2-(se... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.122	2.14 ng/ul	121922	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-(But-1-en-1-yl)-2-(sec-but...	176	C8H16S2	110690-22-7	50		
2	(Z)-1-(But-2-en-1-yl)-2-(sec-but...	176	C8H16S2	110690-23-8	50		
3	(E)-1-(But-2-en-1-yl)-2-(sec-but...	176	C8H16S2	110690-24-9	45		
4	(Z)-1-(But-1-en-1-yl)-2-(sec-but...	176	C8H16S2	110690-21-6	45		
5	Phenylalanine	165	C9H11NO2	000063-91-2	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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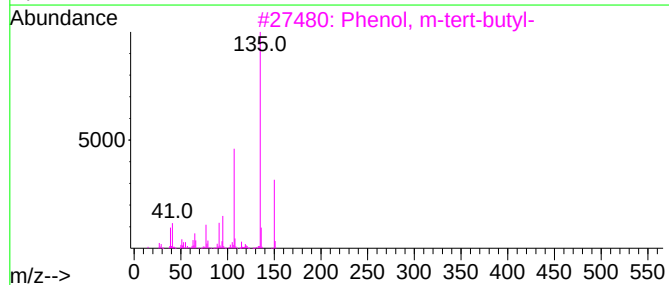
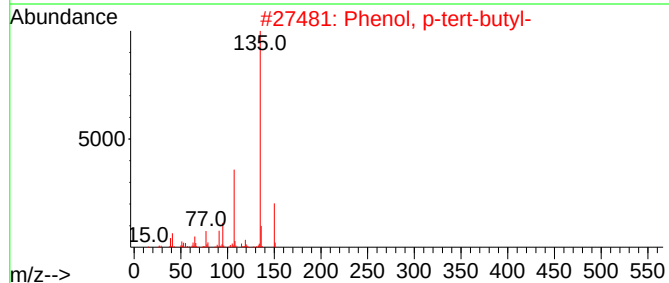
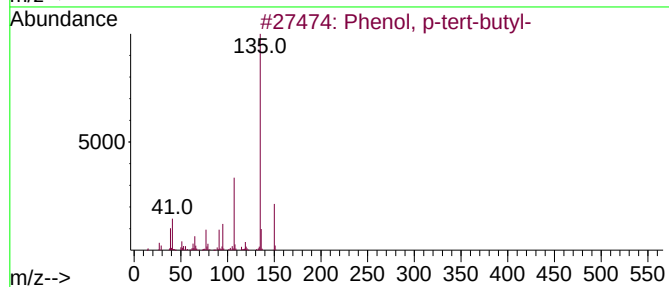
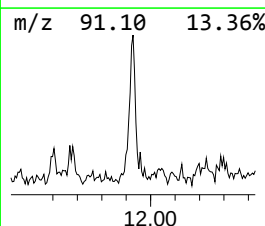
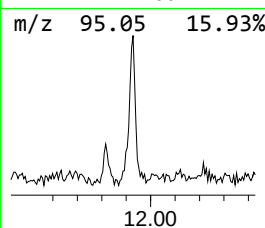
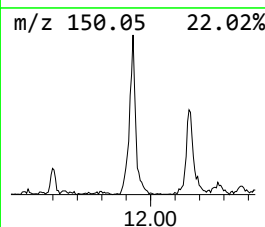
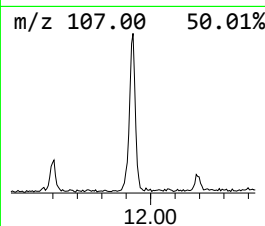
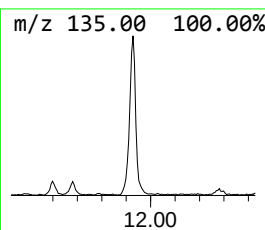
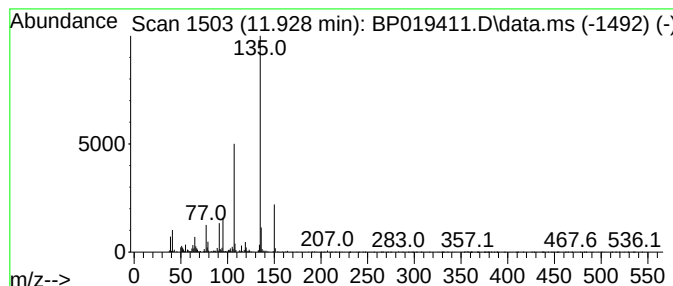
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.928	2.94 ng/ul	166911	Naphthalene-d8	10.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87
5			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

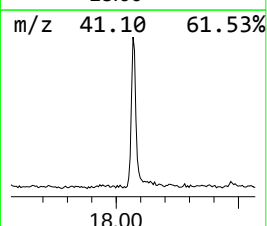
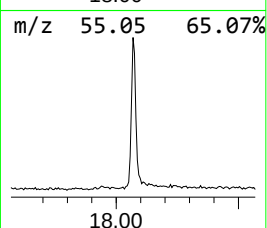
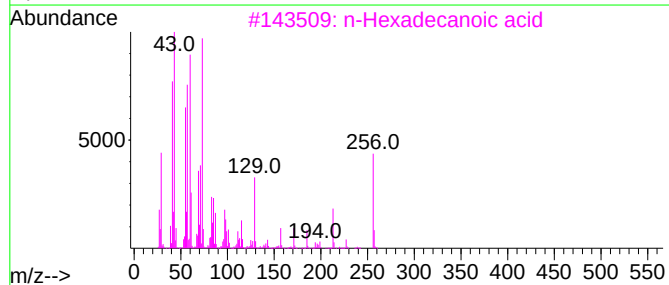
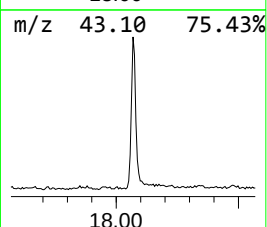
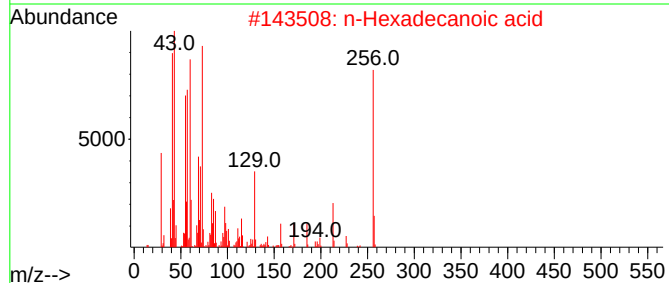
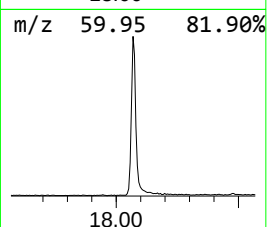
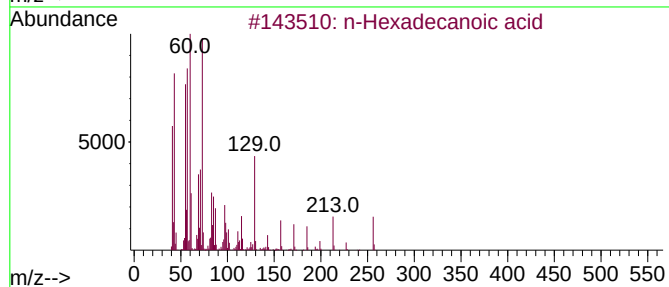
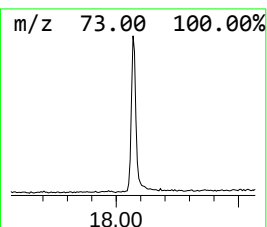
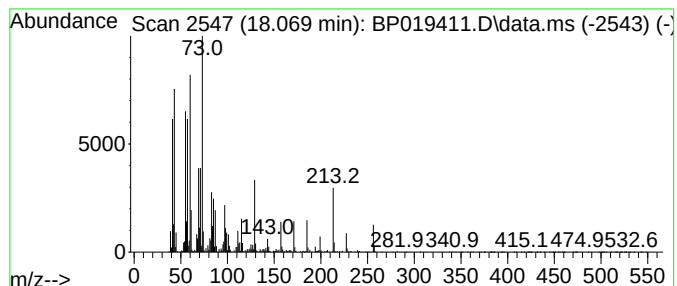
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.069	10.33 ng/ul	880329	Phenanthrene-d10	17.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5	Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

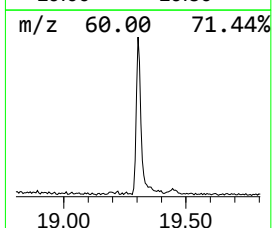
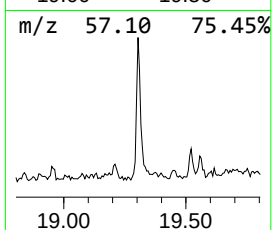
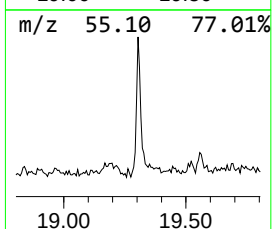
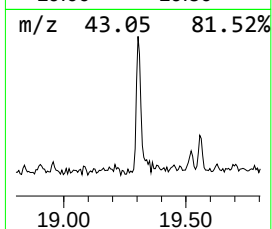
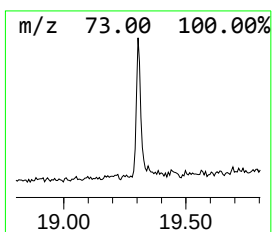
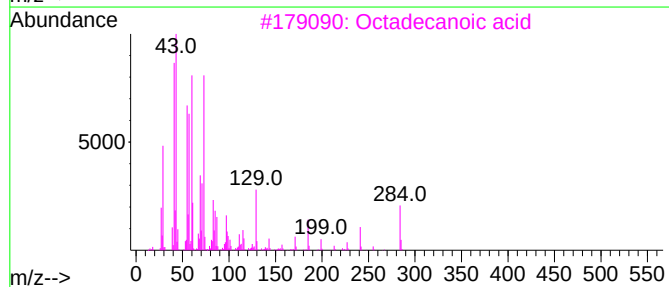
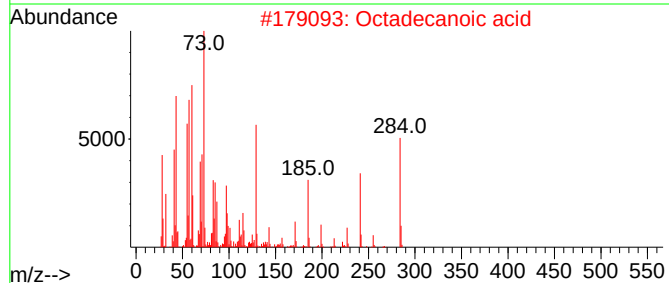
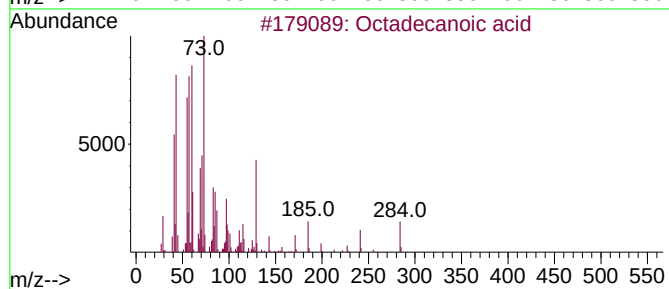
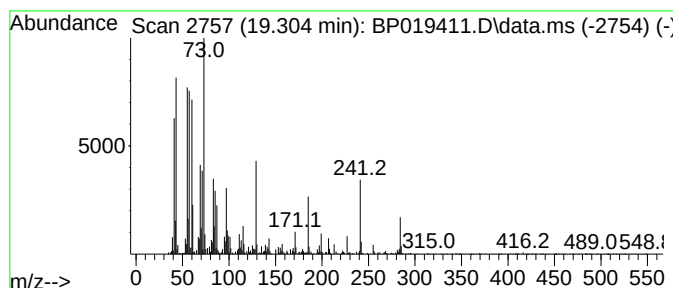
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.304	4.16 ng/ul	312850	Chrysene-d12	21.275	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	98
3	Octadecanoic acid	284	C18H36O2	000057-11-4	98
4	Octadecanoic acid	284	C18H36O2	000057-11-4	97
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

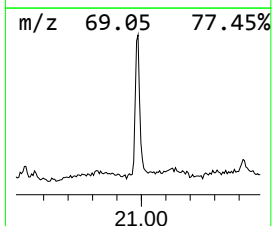
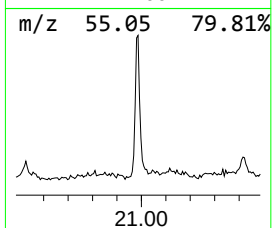
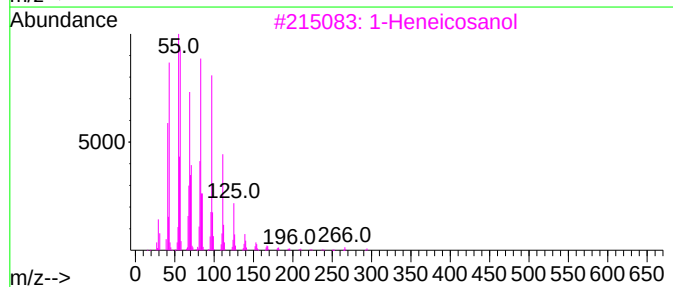
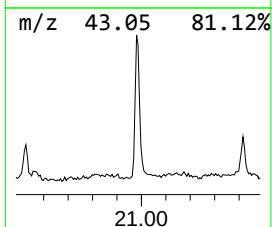
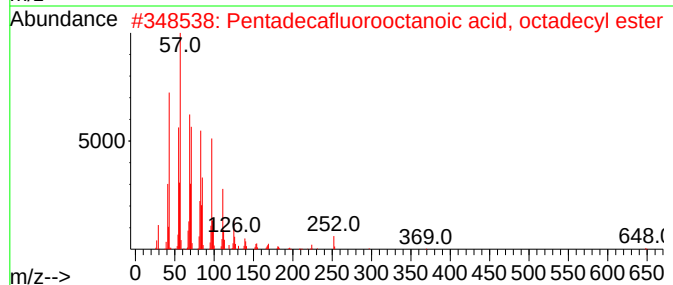
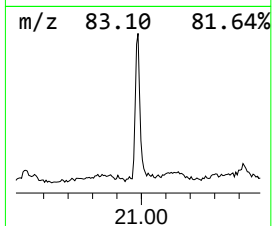
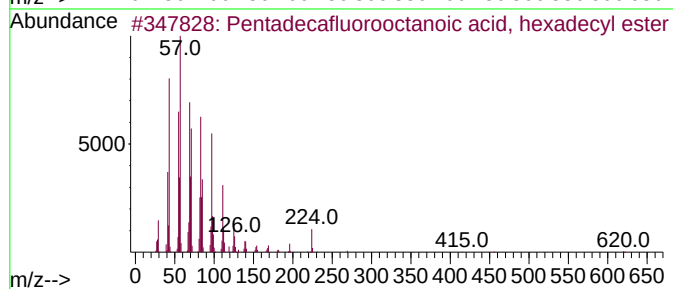
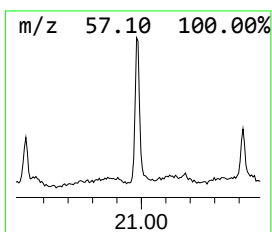
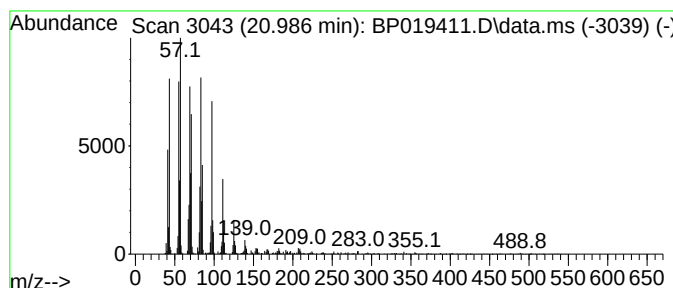
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Pentadecafluorooctanoic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.986	8.24 ng/ul	620345	Chrysene-d12	21.275	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	91
5	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012224\
Data File : BP019411.D
Acq On : 22 Jan 2024 13:34
Operator : MA/JU
Sample : P1142-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.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...	7.658	2.5	ng/ul	105347	1	7.769	830662	20.0
Benzene, 1,2-di...	8.122	2.5	ng/ul	102167	1	7.769	830662	20.0
Benzene, 1,2,4-...	10.428	7.9	ng/ul	447121	2	10.563	1136820	20.0
(E)-1-(But-1-en...	11.122	2.1	ng/ul	121922	2	10.563	1136820	20.0
Phenol, p-tert-...	11.928	2.9	ng/ul	166911	2	10.563	1136820	20.0
n-Hexadecanoic ...	18.069	10.3	ng/ul	880329	4	17.169	1704310	20.0
Octadecanoic acid	19.304	4.2	ng/ul	312850	5	21.275	1505290	20.0
Pentadecafluoro...	20.986	8.2	ng/ul	620345	5	21.275	1505290	20.0