

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
 Data File : BP016233.D
 Acq On : 14 Jul 2023 19:05
 Operator : MA/JU
 Sample : 03414-19
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4634

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Signal : TIC: BP016233.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	13	17	27	rVB	38470	68592	0.84%	0.084%
2	3.740	92	94	118	rBV	346959	830825	10.19%	1.017%
3	4.046	144	146	152	rVB3	10191	15788	0.19%	0.019%
4	7.158	667	675	693	rBV	413553	1333076	16.35%	1.632%
5	7.293	693	698	725	rVV	414372	1048389	12.86%	1.283%
6	7.499	726	733	757	rVV	541906	1231389	15.11%	1.507%
7	7.669	758	762	763	rBV3	7787	9385	0.12%	0.011%
8	7.963	805	812	834	rBV	443660	1122523	13.77%	1.374%
9	8.452	890	895	904	rBV	4938	16155	0.20%	0.020%
10	8.599	914	920	924	rBV7	8245	14802	0.18%	0.018%
11	8.711	932	939	971	rBV2	611615	1870040	22.94%	2.289%
12	9.046	995	996	1004	rVB8	7393	10567	0.13%	0.013%
13	9.158	1006	1015	1043	rBV	524819	1425847	17.49%	1.745%
14	9.499	1067	1073	1081	rVB3	11420	26138	0.32%	0.032%
15	9.663	1097	1101	1106	rBV7	6172	12192	0.15%	0.015%
16	9.816	1122	1127	1130	rBV7	4328	8485	0.10%	0.010%
17	9.893	1130	1140	1165	rBV2	406244	1251982	15.36%	1.532%
18	10.128	1176	1180	1185	rVB2	6445	8253	0.10%	0.010%
19	10.457	1229	1236	1272	rBV	422754	1875453	23.01%	2.295%
20	10.781	1284	1291	1308	rVB	880740	1796023	22.03%	2.198%
21	10.975	1317	1324	1348	rBV2	411832	1689682	20.73%	2.068%
22	11.404	1394	1397	1400	rBV5	7876	11342	0.14%	0.014%
23	11.446	1400	1404	1410	rVB5	14370	28938	0.36%	0.035%
24	11.681	1441	1444	1449	rVB7	6608	8519	0.10%	0.010%
25	11.763	1453	1458	1463	rBV	58336	97727	1.20%	0.120%
26	12.028	1499	1503	1507	rVB6	9010	13181	0.16%	0.016%
27	12.363	1555	1560	1568	rVB6	11232	34426	0.42%	0.042%
28	12.446	1569	1574	1576	rBV4	10796	17934	0.22%	0.022%
29	12.610	1598	1602	1603	rBV4	6975	9743	0.12%	0.012%
30	12.799	1630	1634	1640	rBV8	10547	21258	0.26%	0.026%
31	12.863	1641	1645	1654	rVB4	16253	31540	0.39%	0.039%
32	13.110	1683	1687	1693	rVB	108385	147785	1.81%	0.181%
33	13.363	1726	1730	1731	rBV4	8213	11435	0.14%	0.014%
34	13.381	1731	1733	1737	rVB5	11813	13272	0.16%	0.016%
35	13.422	1737	1740	1745	rBV4	8508	12382	0.15%	0.015%
36	13.487	1745	1751	1762	rBV	997348	1635031	20.06%	2.001%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
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37	13.687	1783	1785	1788	rBV4	6042	8596	0.11%	0.011%
38	13.793	1798	1803	1807	rBV6	18642	34132	0.42%	0.042%
39	13.910	1819	1823	1824	rBV4	7415	8730	0.11%	0.011%
40	14.028	1835	1843	1857	rBV	1762878	3289008	40.35%	4.026%
41	14.310	1885	1891	1904	rBV	2328568	3773204	46.29%	4.618%
42	14.410	1905	1908	1914	rVB4	56836	87983	1.08%	0.108%
43	14.546	1927	1931	1935	rBV5	26195	39199	0.48%	0.048%
44	14.610	1937	1942	1951	rVV	1621641	2508054	30.77%	3.070%
45	14.734	1957	1963	1975	rBV	1304607	2202065	27.02%	2.695%
46	14.928	1990	1996	2002	rBV	247069	659326	8.09%	0.807%
47	15.304	2056	2060	2068	rVB7	25183	59008	0.72%	0.072%
48	15.375	2068	2072	2081	rBV10	26514	69389	0.85%	0.085%
49	15.616	2104	2113	2128	rBV	2641653	4953780	60.78%	6.063%
50	15.798	2139	2144	2148	rBV	404977	813274	9.98%	0.995%
51	15.863	2150	2155	2173	rVB	2246190	3573945	43.85%	4.374%
52	16.045	2184	2186	2190	rVB5	13521	13020	0.16%	0.016%
53	16.310	2226	2231	2236	rBV	242764	359067	4.41%	0.439%
54	16.481	2256	2260	2268	rBV	83041	122995	1.51%	0.151%
55	16.598	2275	2280	2285	rBV	387204	573189	7.03%	0.702%
56	16.887	2324	2329	2336	rBV	382505	572868	7.03%	0.701%
57	17.139	2367	2372	2381	rBV	160767	245290	3.01%	0.300%
58	17.239	2385	2389	2391	rVV3	63721	92308	1.13%	0.113%
59	17.275	2391	2395	2402	rVB2	185338	275384	3.38%	0.337%
60	17.375	2407	2412	2418	rBV	2267564	3584484	43.98%	4.387%
61	17.481	2424	2430	2439	rBV2	3049426	5420627	66.50%	6.635%
62	17.545	2439	2441	2451	rVB	203624	289912	3.56%	0.355%
63	17.745	2472	2475	2480	rBV2	56375	76014	0.93%	0.093%
64	17.828	2486	2489	2492	rBV	109044	140289	1.72%	0.172%
65	17.963	2509	2512	2519	rBV2	124986	250928	3.08%	0.307%
66	18.069	2527	2530	2534	rBV3	89355	112487	1.38%	0.138%
67	18.151	2541	2544	2548	rVB5	60259	77895	0.96%	0.095%
68	18.239	2554	2559	2569	rBV	1007137	1606528	19.71%	1.966%
69	18.416	2586	2589	2593	rBV2	105147	150771	1.85%	0.185%
70	18.481	2597	2600	2603	rVV3	122002	163074	2.00%	0.200%
71	18.516	2603	2606	2611	rVB2	159886	215701	2.65%	0.264%
72	18.786	2649	2652	2655	rVB4	66924	69964	0.86%	0.086%
73	19.139	2709	2712	2716	rVB2	147540	158063	1.94%	0.193%
74	19.392	2752	2755	2761	rVB	197411	265563	3.26%	0.325%
75	19.463	2763	2767	2777	rBV	731962	1121469	13.76%	1.373%
76	19.716	2804	2810	2826	rBV	4392862	7455218	91.46%	9.125%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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

77	20.186	2888	2890	2897	rVB	103221	103735	1.27%	0.127%
78	21.333	3082	3085	3092	rVB2	393668	509104	6.25%	0.623%
79	21.475	3104	3109	3121	rBV	2625738	4430517	54.36%	5.423%
80	22.033	3201	3204	3208	rVB	213992	254395	3.12%	0.311%
81	23.110	3383	3387	3395	rVB	395643	621141	7.62%	0.760%
82	23.798	3498	3504	3523	rVB2	3600934	8150919	100.00%	9.976%
83	23.969	3526	3533	3546	rVB	1735217	4378909	53.72%	5.360%

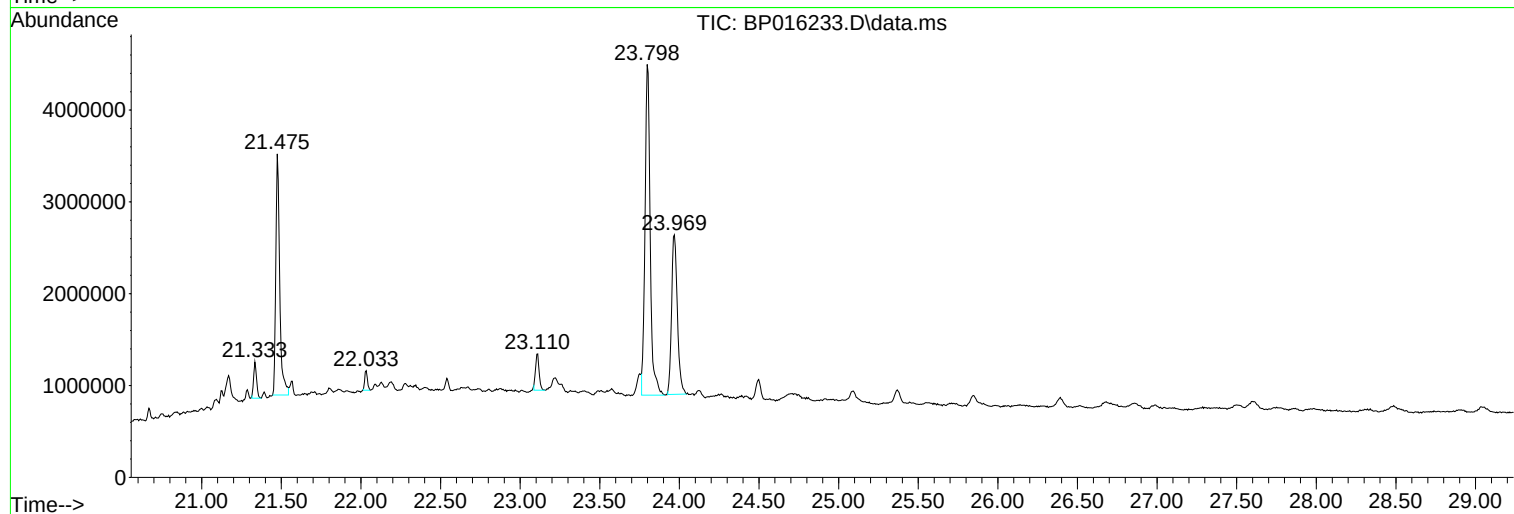
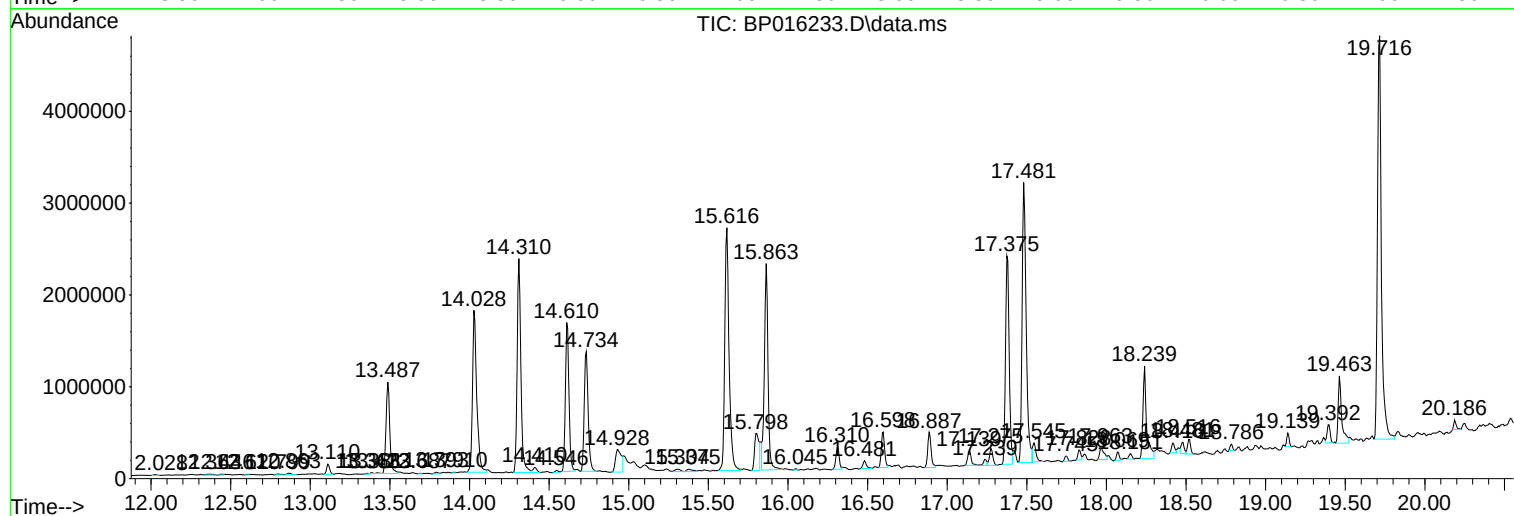
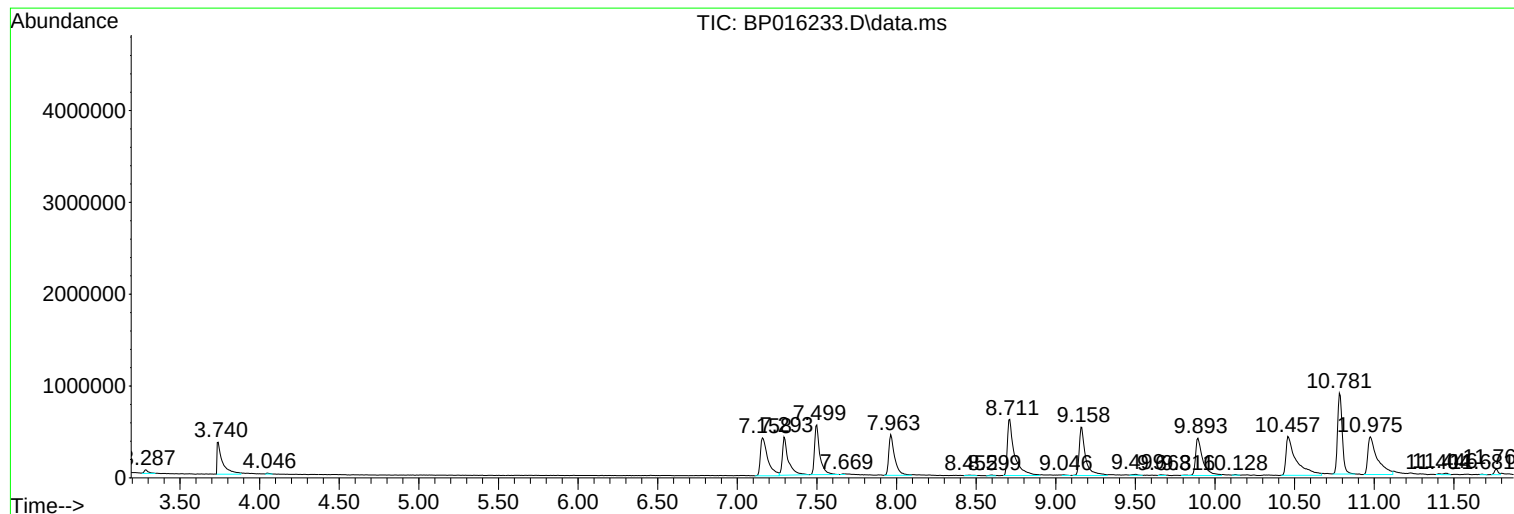
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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



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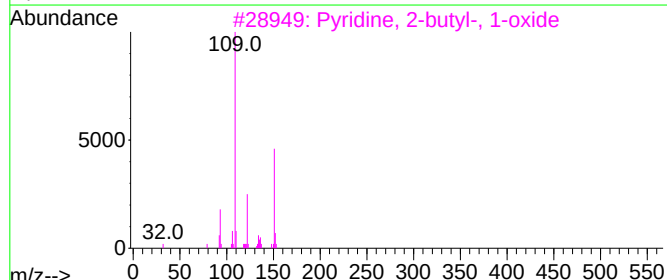
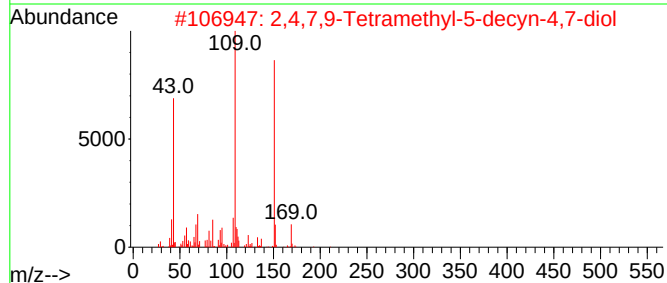
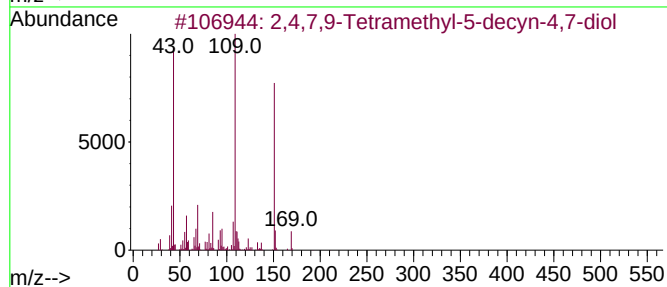
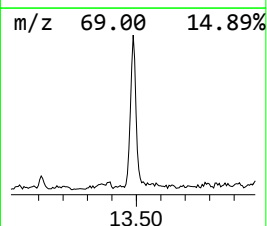
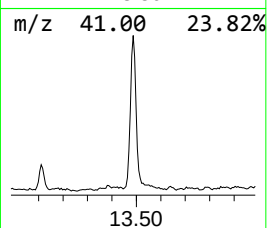
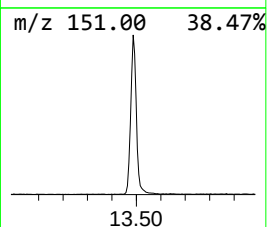
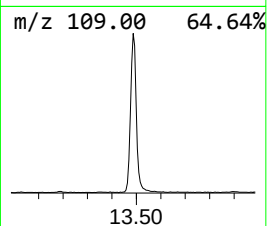
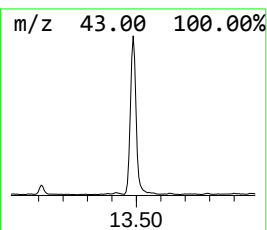
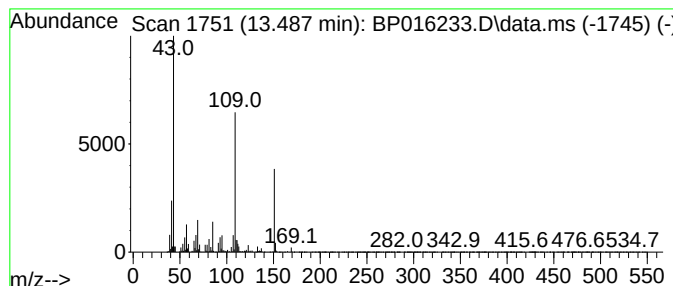
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.487	13.04 ng/ul	1635030	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	74		
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
4	Acetaminophen	151	C8H9NO2	000103-90-2	53		
5	Metacetamol	151	C8H9NO2	000621-42-1	53		



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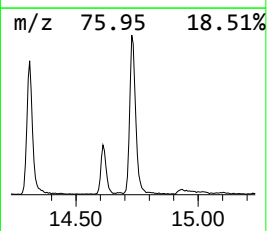
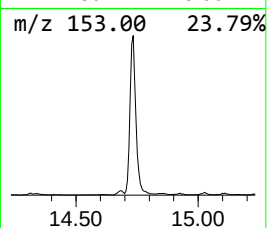
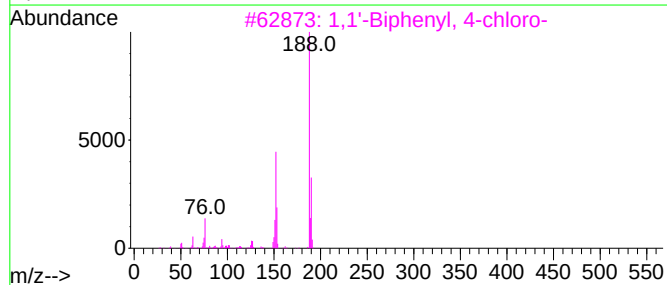
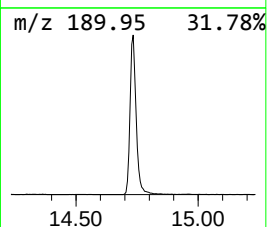
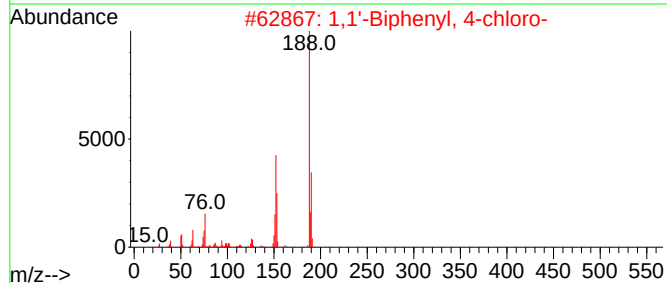
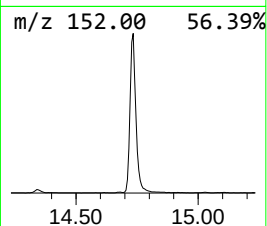
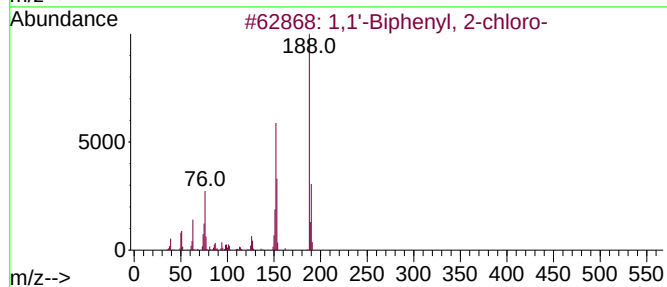
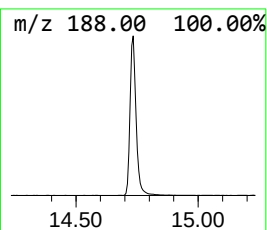
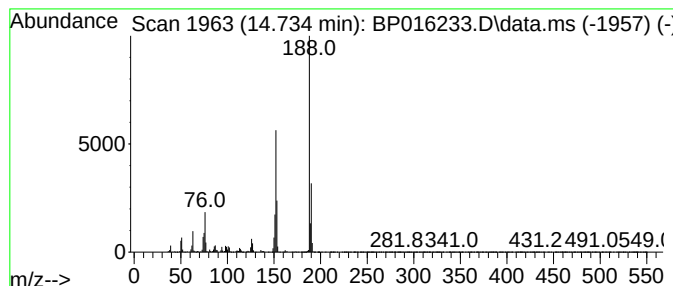
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.734	17.56 ng/ul	2202070	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-chloro-			188	C12H9Cl	002051-60-7	98
2	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	97
3	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	95
4	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	95
5	3-Chlorobiphenyl			188	C12H9Cl	002051-61-8	94



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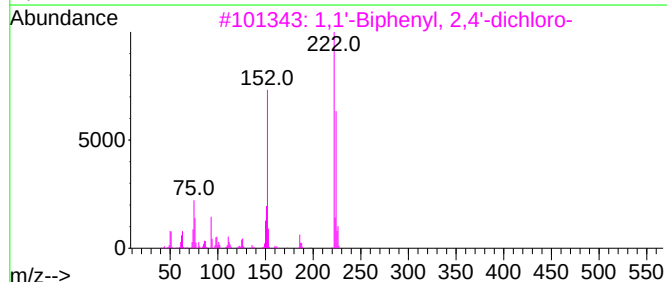
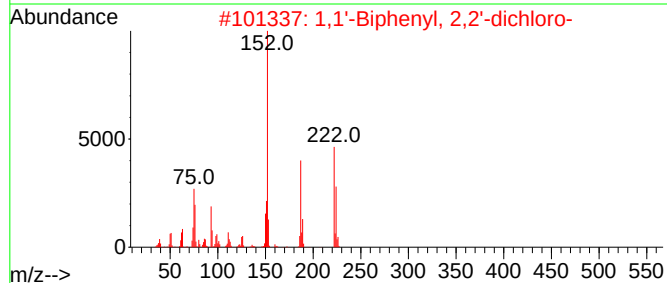
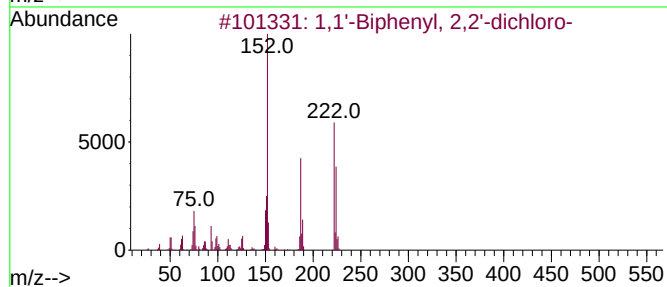
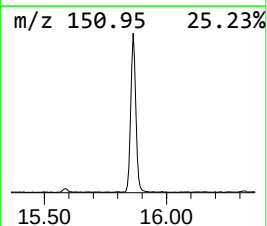
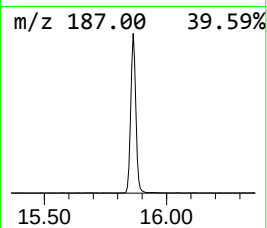
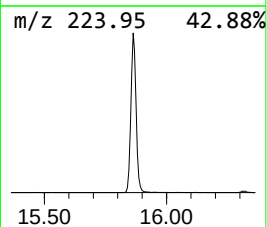
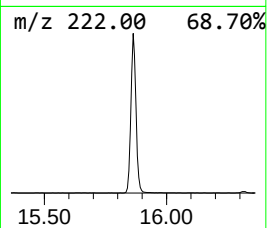
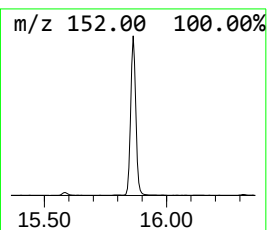
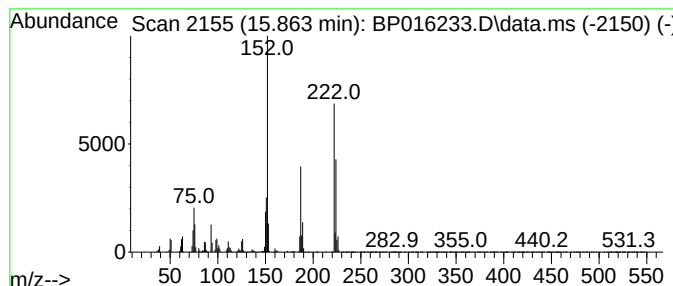
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.863	28.50 ng/ul	3573950	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99		
2	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98		
3	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	98		
4	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98		
5	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	98		



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A4634

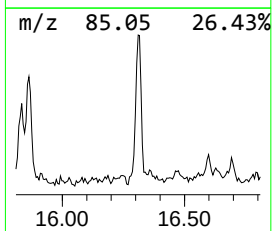
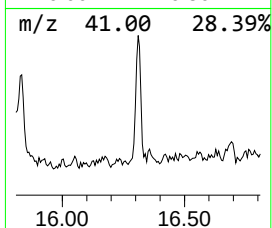
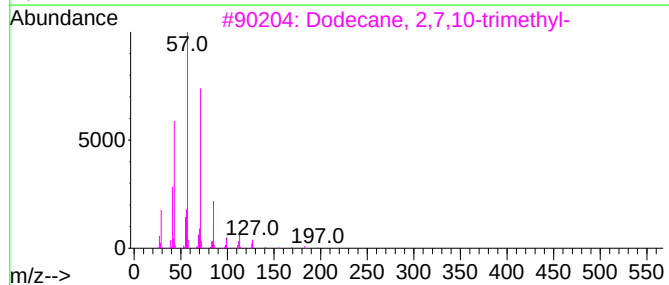
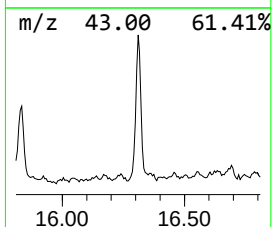
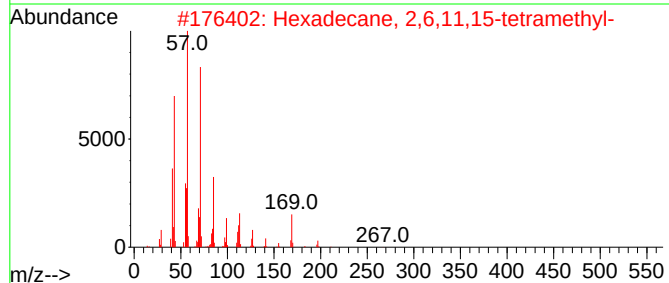
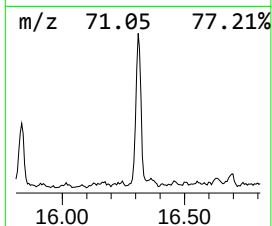
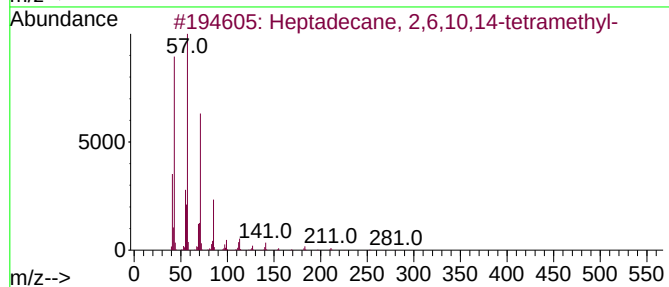
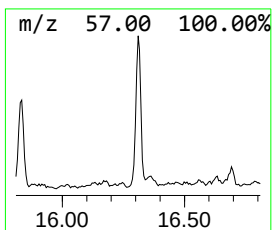
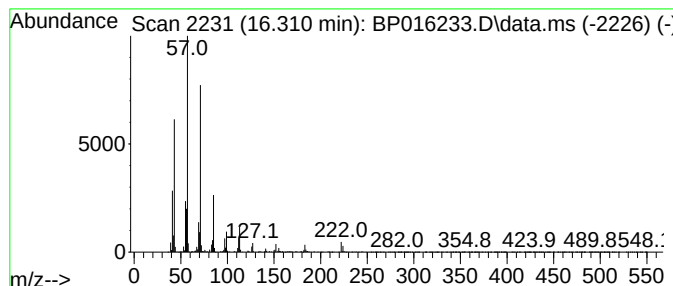
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.310	2.00 ng/ul	359067	Phenanthrene-d10	17.375

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	91
2		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	87
3		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	87
4		Decane, 2-methyl-	156	C11H24	006975-98-0	87
5		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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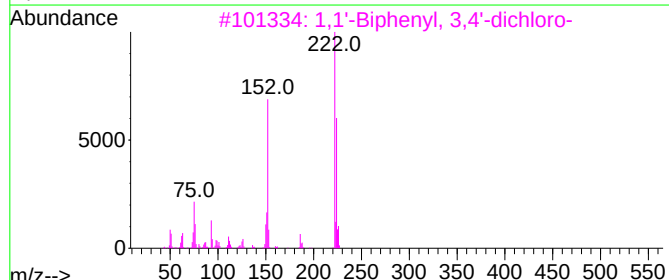
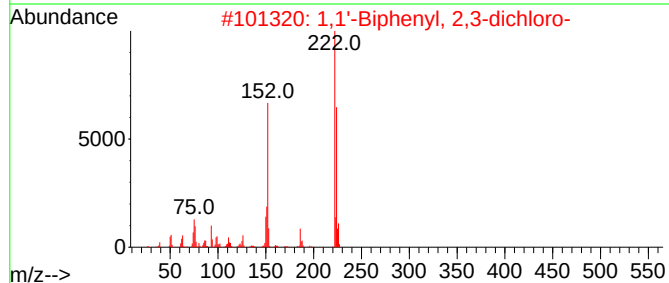
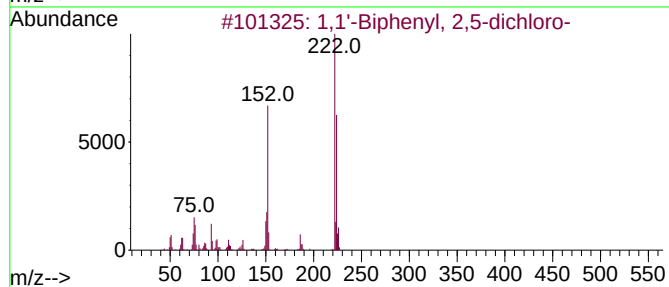
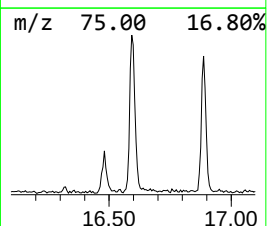
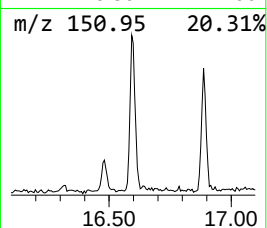
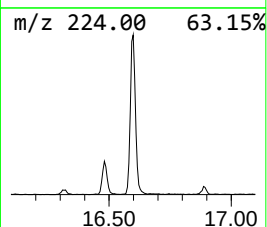
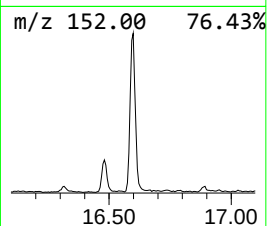
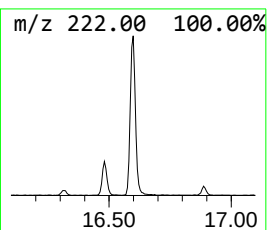
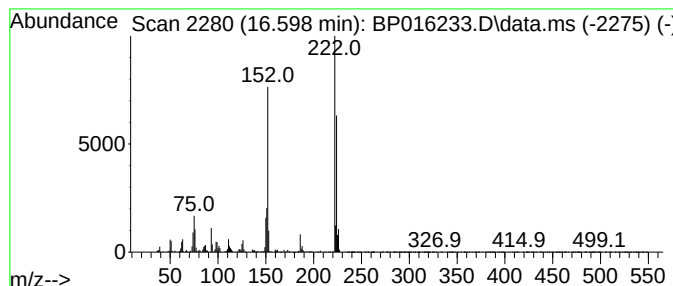
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,5-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	3.20 ng/ul	573189	Phenanthrene-d10	17.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	99		
2	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99		
3	1,1'-Biphenyl, 3,4'-dichloro-	222	C12H8Cl2	002974-90-5	99		
4	1,1'-Biphenyl, 3,4-dichloro-	222	C12H8Cl2	002974-92-7	99		
5	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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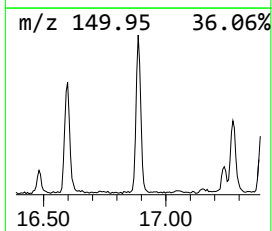
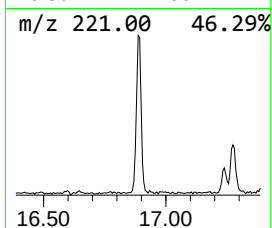
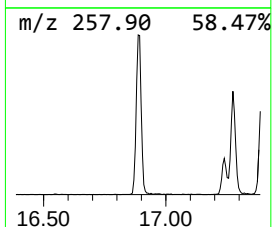
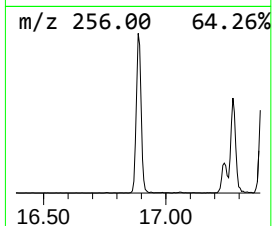
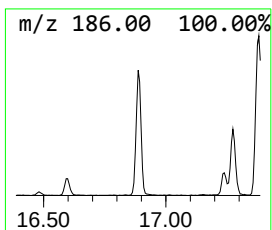
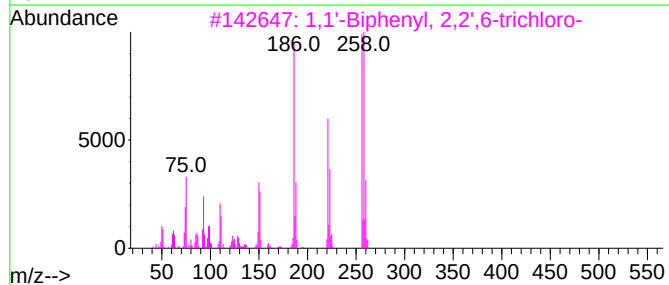
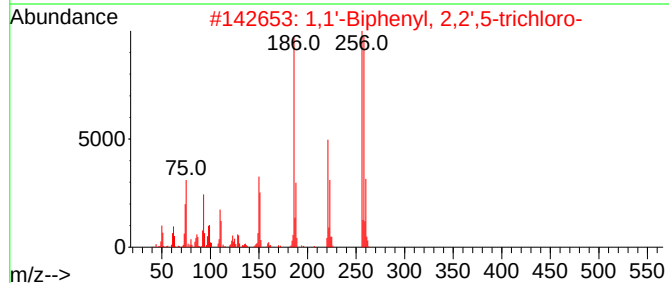
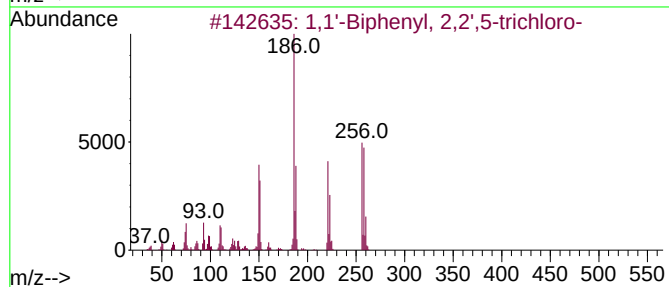
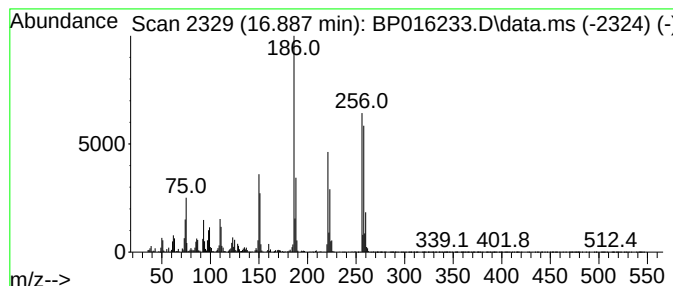
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.887	3.20 ng/ul	572868	Phenanthrene-d10	17.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
3	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	98		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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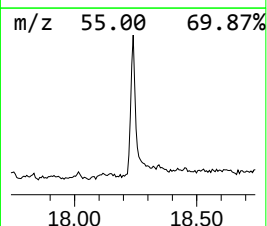
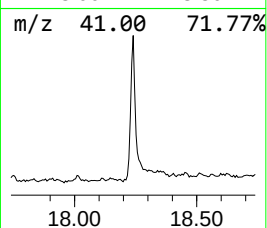
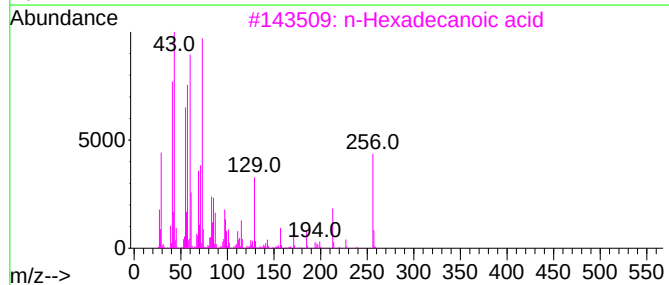
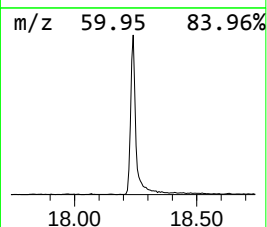
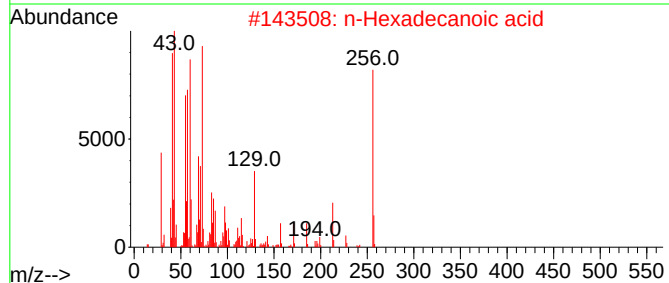
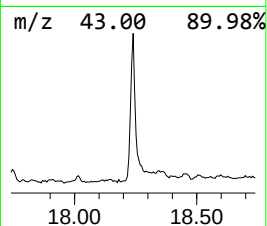
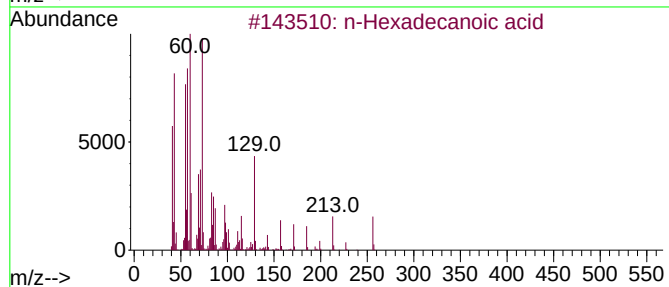
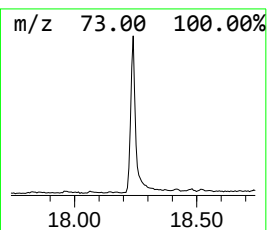
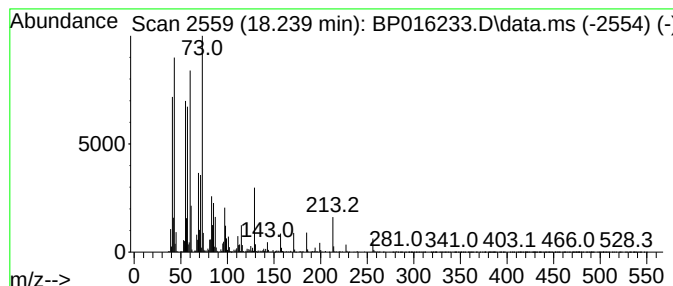
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	8.96 ng/ul	1606530	Phenanthrene-d10	17.375

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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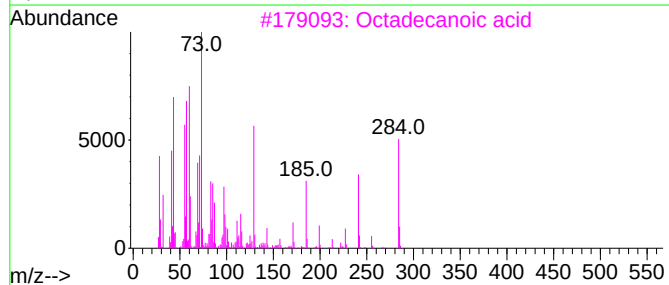
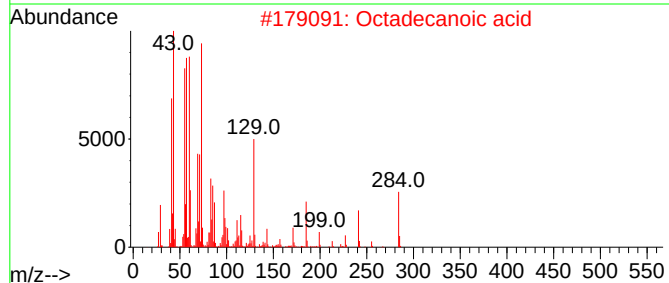
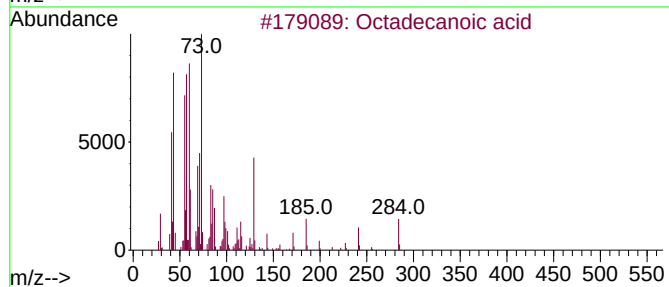
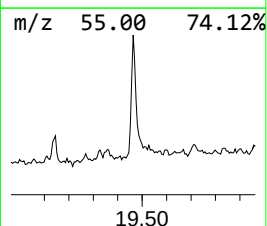
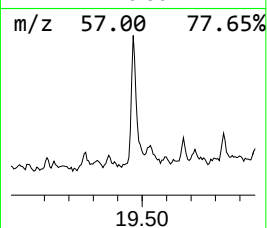
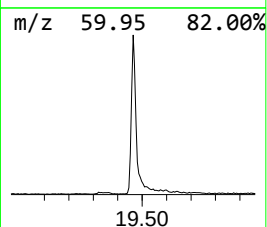
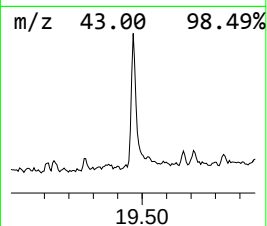
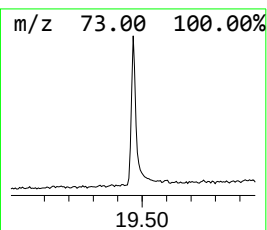
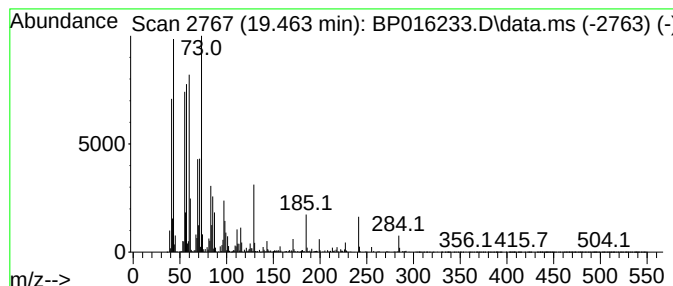
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	5.06 ng/ul	1121470	Chrysene-d12	21.474

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	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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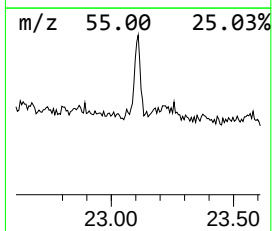
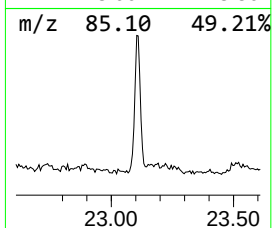
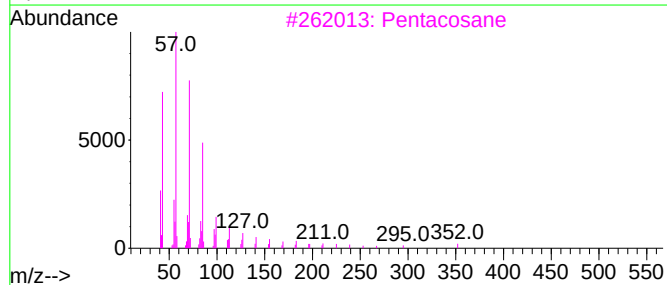
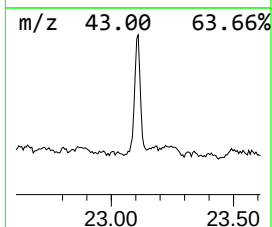
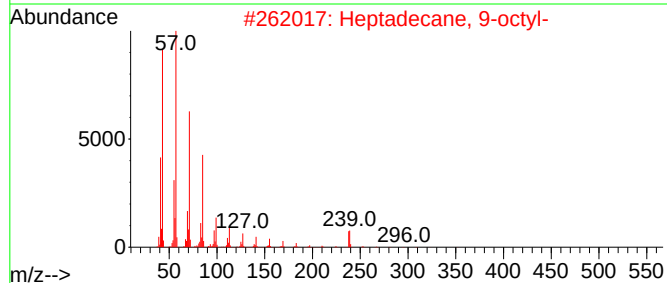
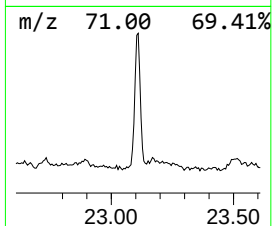
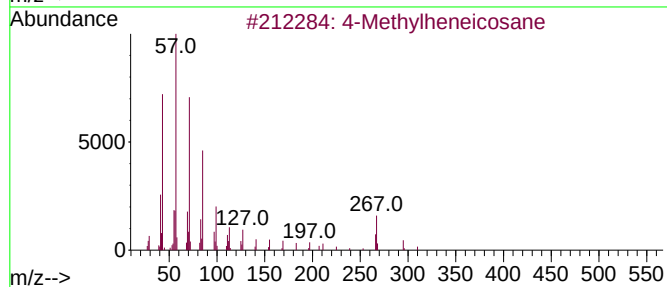
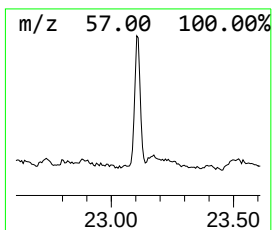
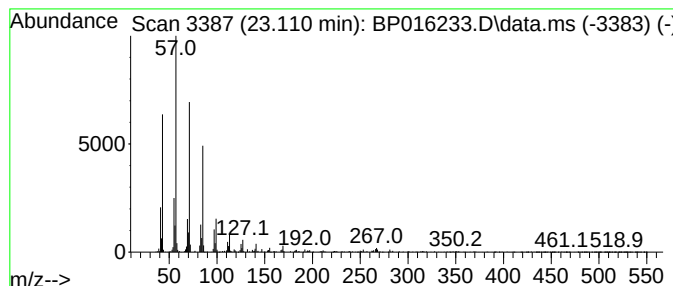
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.110	2.84 ng/ul	621141	Perylene-d12	23.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylheneicosane	310	C22H46	025117-29-7	93
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			Pentacosane	352	C25H52	000629-99-2	93
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5			Hexadecane	226	C16H34	000544-76-3	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016233.D
Acq On : 14 Jul 2023 19:05
Operator : MA/JU
Sample : 03414-19
Misc :
ALS Vial : 18 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,4,7,9-Tetrame...	13.487	13.0	ng/ul	1635030	3	14.610	2508050	20.0
1,1'-Biphenyl, ...	14.734	17.6	ng/ul	2202070	3	14.610	2508050	20.0
1,1'-Biphenyl, ...	15.863	28.5	ng/ul	3573950	3	14.610	2508050	20.0
(DEL) Alkane: S...	16.310	2.0	ng/ul	359067	4	17.375	3584480	20.0
1,1'-Biphenyl, ...	16.598	3.2	ng/ul	573189	4	17.375	3584480	20.0
1,1'-Biphenyl, ...	16.887	3.2	ng/ul	572868	4	17.375	3584480	20.0
n-Hexadecanoic ...	18.239	9.0	ng/ul	1606530	4	17.375	3584480	20.0
Octadecanoic acid	19.463	5.1	ng/ul	1121470	5	21.474	4430520	20.0
(DEL) Alkane: S...	23.110	2.8	ng/ul	621141	6	23.968	4378910	20.0