

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
 Data File : BM036478.D
 Acq On : 01 Sep 2022 18:12
 Operator : CG/JU
 Sample : N4458-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COKW0

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_M\Methods\SFAM-EPA-BM081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036478.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	20	24	34	rVB	50189	73949	0.99%	0.096%
2	3.563	95	98	114	rBV	161849	301763	4.04%	0.391%
3	3.781	130	135	141	rVB	24711	40070	0.54%	0.052%
4	3.875	147	151	156	rBV	23443	34588	0.46%	0.045%
5	5.022	339	346	356	rBV	330706	505444	6.77%	0.654%
6	6.910	661	667	679	rBV	159827	286568	3.84%	0.371%
7	7.069	684	694	706	rBV	753085	1154015	15.45%	1.494%
8	7.257	719	726	737	rBV	680646	1096562	14.68%	1.419%
9	7.333	737	739	751	rVV9	9756	23183	0.31%	0.030%
10	7.439	753	757	766	rVB5	5819	11688	0.16%	0.015%
11	7.610	781	786	793	rBV	90290	146830	1.97%	0.190%
12	7.728	798	806	809	rBV	653128	1048363	14.04%	1.357%
13	7.763	809	812	822	rVB	336668	497692	6.66%	0.644%
14	8.075	857	865	876	rBV	2123763	3290498	44.05%	4.259%
15	8.451	923	929	942	rBV	517598	834221	11.17%	1.080%
16	8.898	997	1005	1008	rBV	902262	1470024	19.68%	1.903%
17	8.939	1008	1012	1023	rVB	1042947	1649832	22.09%	2.135%
18	9.057	1030	1032	1040	rVB6	5376	8061	0.11%	0.010%
19	9.286	1066	1071	1075	rBV	15534	23489	0.31%	0.030%
20	9.357	1078	1083	1088	rVB8	8337	13924	0.19%	0.018%
21	9.539	1110	1114	1118	rBV5	6538	11244	0.15%	0.015%
22	9.616	1120	1127	1145	rVV	733564	1240800	16.61%	1.606%
23	10.151	1211	1218	1229	rBV	974889	1637674	21.93%	2.120%
24	10.222	1229	1230	1238	rVV7	12342	18544	0.25%	0.024%
25	10.316	1240	1246	1251	rBV10	6047	10541	0.14%	0.014%
26	10.386	1251	1258	1266	rBV	982058	1546196	20.70%	2.001%
27	10.463	1266	1271	1274	rVV5	10481	23065	0.31%	0.030%
28	10.522	1274	1281	1289	rVB	958093	1599632	21.42%	2.070%
29	10.674	1299	1307	1319	rBV	488732	921855	12.34%	1.193%
30	10.910	1337	1347	1359	rVB	512297	856264	11.46%	1.108%
31	11.139	1379	1386	1403	rVB	704905	1173219	15.71%	1.519%
32	11.357	1415	1423	1434	rBV	143696	294447	3.94%	0.381%
33	11.545	1450	1455	1459	rBV6	5007	10621	0.14%	0.014%
34	11.780	1490	1495	1500	rBV	32112	50677	0.68%	0.066%
35	11.839	1500	1505	1516	rVV2	61139	119605	1.60%	0.155%
36	11.951	1518	1524	1530	rVB	25755	38866	0.52%	0.050%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
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37	12.121	1548	1553	1559	rVB2	20206	31574	0.42%	0.041%
38	12.451	1602	1609	1615	rBV7	9754	17224	0.23%	0.022%
39	12.527	1615	1622	1625	rVV	79128	130358	1.75%	0.169%
40	12.563	1625	1628	1637	rVB2	48912	76353	1.02%	0.099%
41	12.768	1659	1663	1666	rVB4	7874	10154	0.14%	0.013%
42	12.815	1666	1671	1676	rBV3	27126	49364	0.66%	0.064%
43	12.910	1682	1687	1692	rVB2	20650	28594	0.38%	0.037%
44	12.974	1692	1698	1705	rVB5	25016	48907	0.65%	0.063%
45	13.174	1725	1732	1735	rBV	295741	499750	6.69%	0.647%
46	13.298	1749	1753	1757	rBV7	7120	13474	0.18%	0.017%
47	13.351	1757	1762	1767	rVV6	25579	41491	0.56%	0.054%
48	13.421	1767	1774	1786	rVB2	39815	91380	1.22%	0.118%
49	13.721	1821	1825	1830	rBV5	9183	18289	0.24%	0.024%
50	13.798	1830	1838	1844	rVV	2143931	2976643	39.85%	3.853%
51	13.857	1844	1848	1852	rVV	119422	179842	2.41%	0.233%
52	13.898	1852	1855	1863	rVB2	41555	61578	0.82%	0.080%
53	13.974	1864	1868	1872	rBV	20283	25212	0.34%	0.033%
54	14.027	1872	1877	1878	rBV5	16160	21409	0.29%	0.028%
55	14.068	1878	1884	1896	rVB	2530660	3684048	49.32%	4.768%
56	14.204	1902	1907	1914	rBV8	17881	35081	0.47%	0.045%
57	14.274	1914	1919	1927	rBV	400776	502131	6.72%	0.650%
58	14.374	1927	1936	1950	rBV2	1633899	2633355	35.26%	3.408%
59	14.609	1971	1976	1987	rBV2	124207	248858	3.33%	0.322%
60	14.692	1987	1990	1993	rBV4	16635	27461	0.37%	0.036%
61	14.774	2001	2004	2008	rBV	45326	63096	0.84%	0.082%
62	14.833	2012	2014	2016	rVV	24278	24629	0.33%	0.032%
63	14.892	2016	2024	2030	rVV4	107759	219885	2.94%	0.285%
64	14.962	2032	2036	2040	rBV2	32555	41224	0.55%	0.053%
65	15.009	2041	2044	2048	rVB6	20322	31474	0.42%	0.041%
66	15.174	2070	2072	2076	rBV2	11147	13971	0.19%	0.018%
67	15.227	2076	2081	2086	rVB	618495	719139	9.63%	0.931%
68	15.286	2086	2091	2094	rBV5	16496	31497	0.42%	0.041%
69	15.327	2094	2098	2099	rBV3	17886	22083	0.30%	0.029%
70	15.368	2099	2105	2112	rBV	3391052	4860198	65.07%	6.291%
71	15.504	2123	2128	2139	rVB	1154075	1526213	20.43%	1.975%
72	15.633	2143	2150	2155	rVV	215600	363902	4.87%	0.471%
73	15.727	2162	2166	2172	rVB4	42110	69348	0.93%	0.090%
74	15.786	2172	2176	2180	rBV3	49288	83521	1.12%	0.108%
75	15.845	2182	2186	2190	rVB2	74255	104479	1.40%	0.135%
76	16.109	2223	2231	2236	rBV	684098	1454331	19.47%	1.882%

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77	16.668	2323	2326	2328	rBV3	65722	88550	1.19%	0.115%
78	16.886	2360	2363	2365	rBV2	137525	167322	2.24%	0.217%
79	16.933	2367	2371	2376	rVV	684902	946994	12.68%	1.226%
80	17.121	2398	2403	2409	rBV	2265983	3241497	43.40%	4.196%
81	17.221	2415	2420	2426	rBV	4501026	6026185	80.68%	7.800%
82	17.545	2471	2475	2479	rBV	465678	695623	9.31%	0.900%
83	18.503	2634	2638	2642	rBV2	851931	1174485	15.72%	1.520%
84	19.521	2807	2811	2818	rVB	5275327	6973859	93.37%	9.027%
85	21.227	3099	3101	3110	rVB2	578471	703299	9.42%	0.910%
86	21.309	3111	3115	3122	rVB	2680488	3341135	44.73%	4.325%
87	23.456	3473	3480	3495	rVB2	3898256	7469239	100.00%	9.668%
88	23.603	3499	3505	3514	rVB2	1732124	3285598	43.99%	4.253%

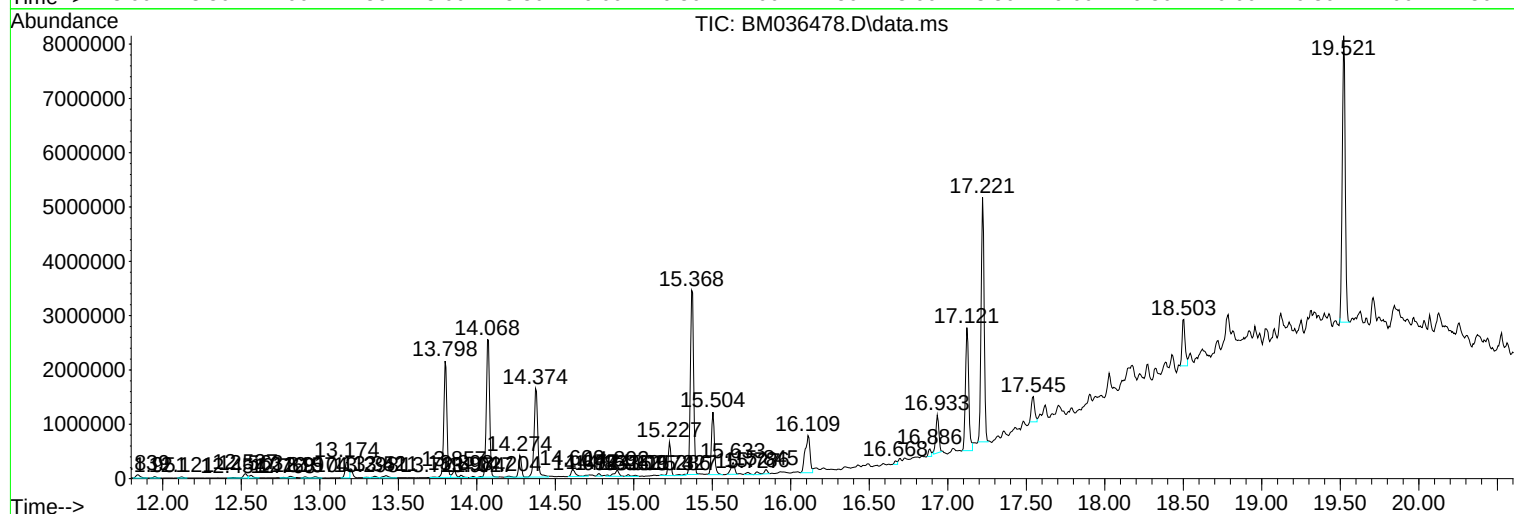
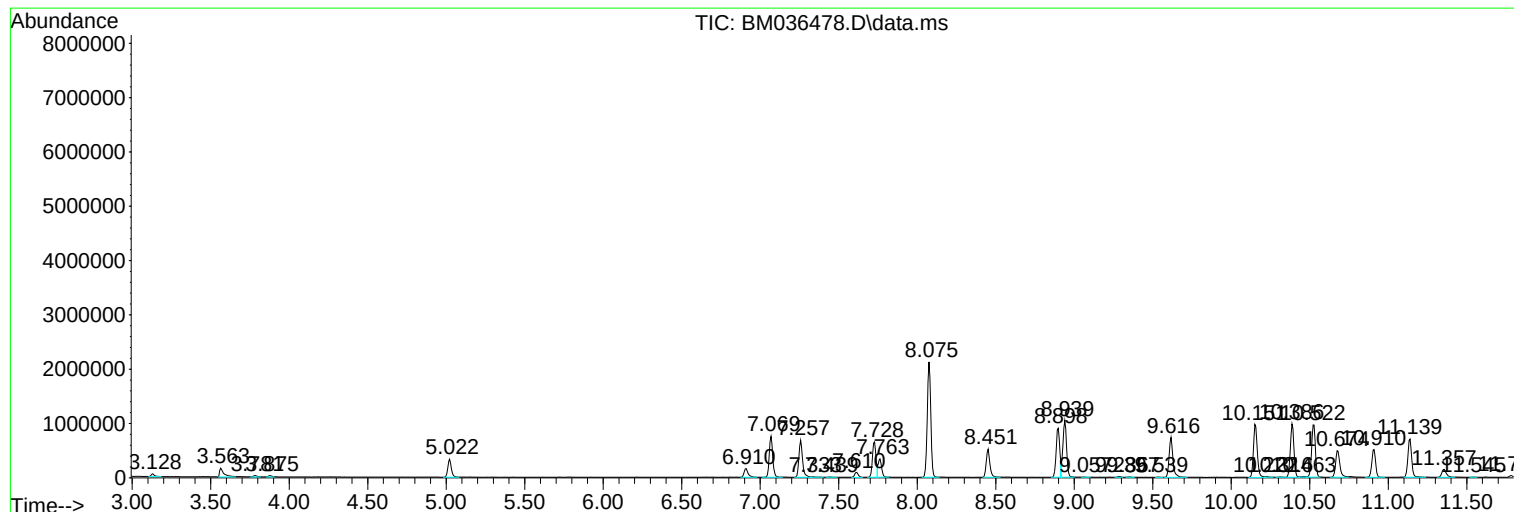
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION

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



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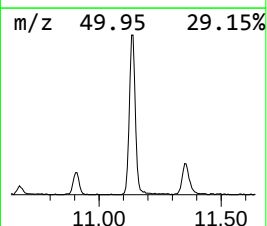
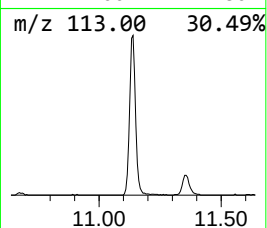
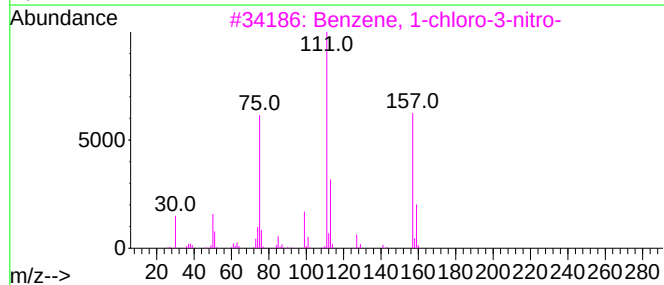
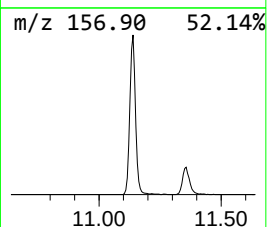
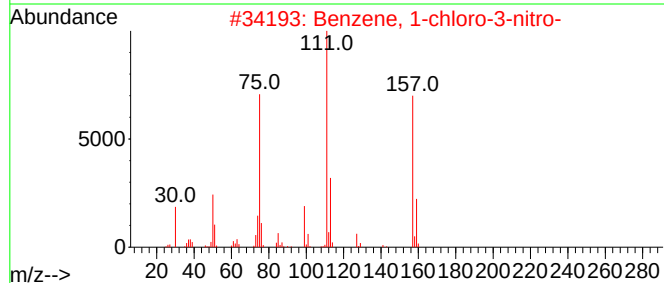
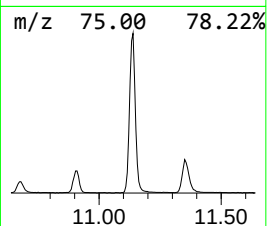
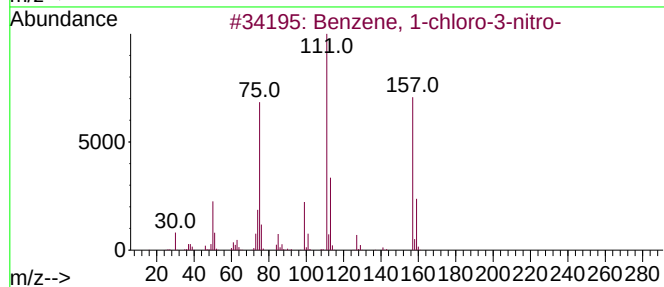
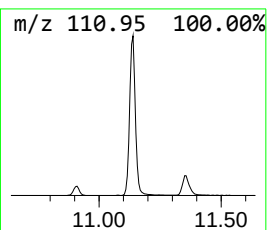
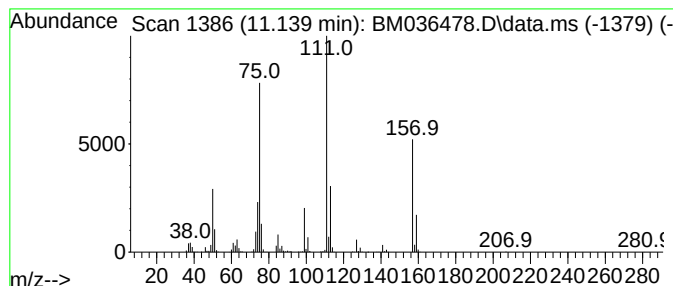
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.139	14.67 ng/ul	1173220	Naphthalene-d8	10.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
2			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	94
3			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	92
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90



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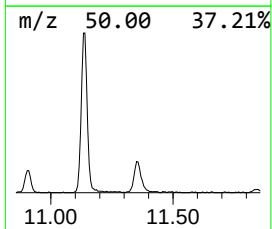
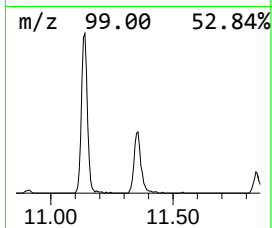
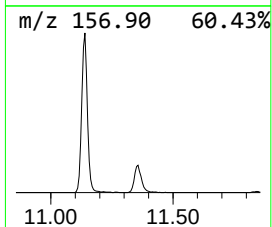
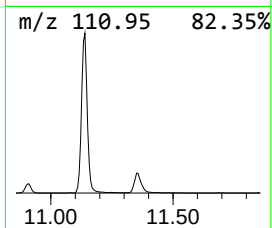
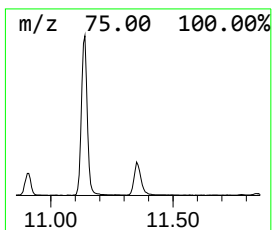
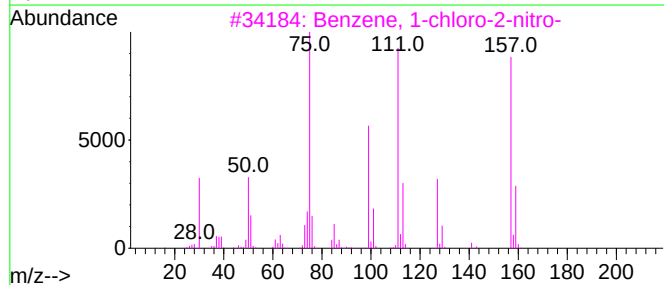
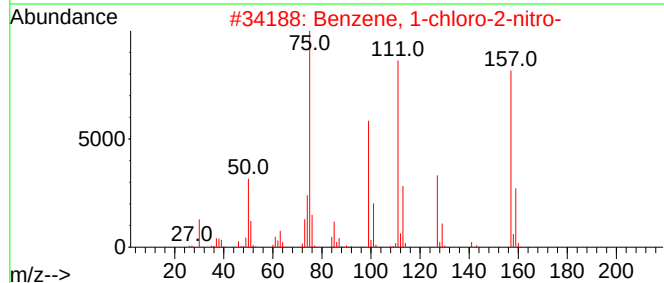
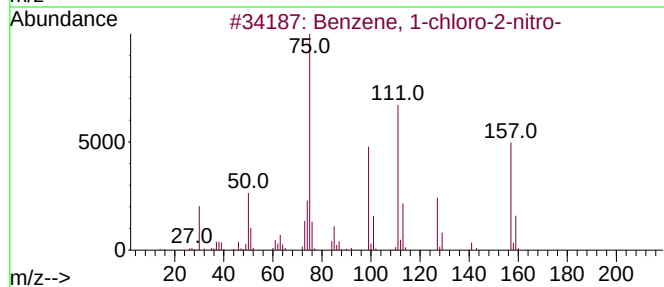
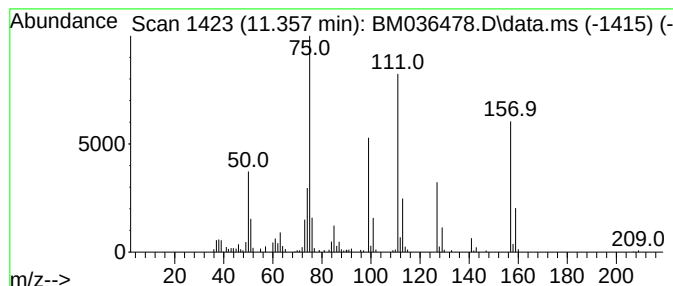
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-chloro-2-nitro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.357	3.68 ng/ul	294447	Naphthalene-d8	10.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
3			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
5			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96



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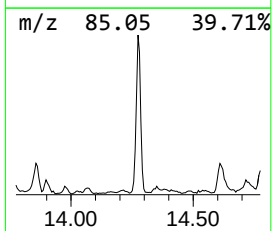
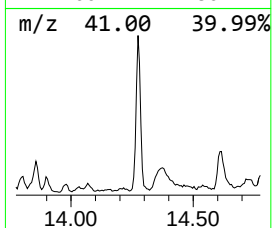
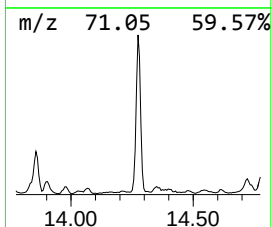
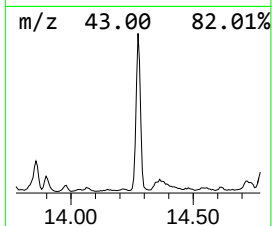
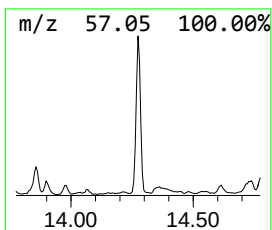
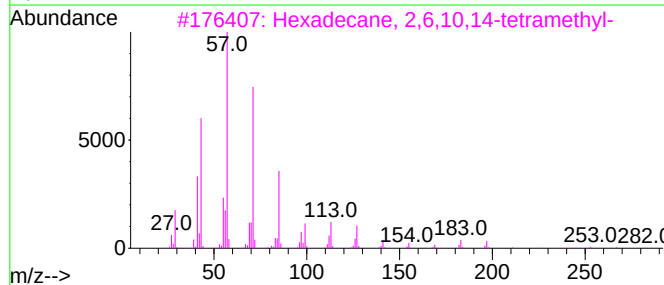
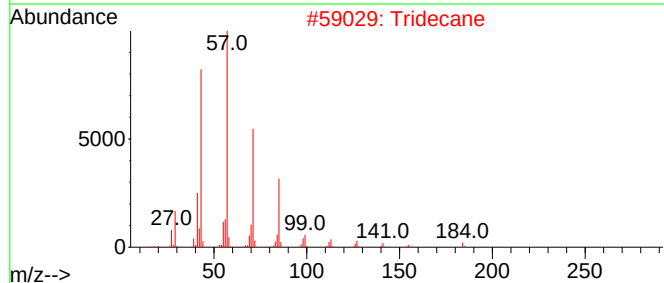
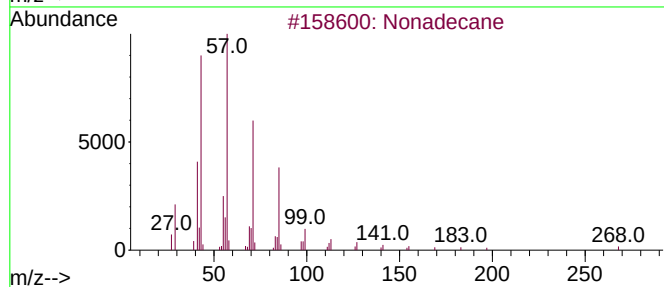
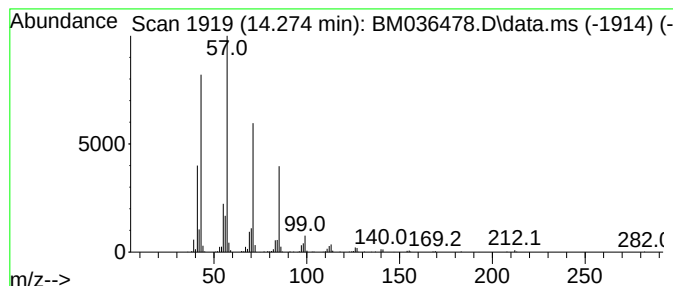
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.274	3.81 ng/ul	502131	Acenaphthene-d10	14.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	90
2	Tridecane	184	C13H28	000629-50-5	86
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86
4	Hexadecane	226	C16H34	000544-76-3	86
5	Tetradecane	198	C14H30	000629-59-4	86



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COKW0

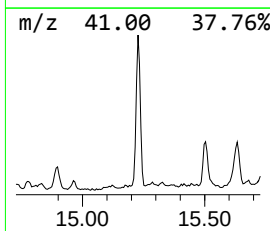
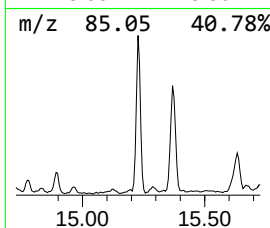
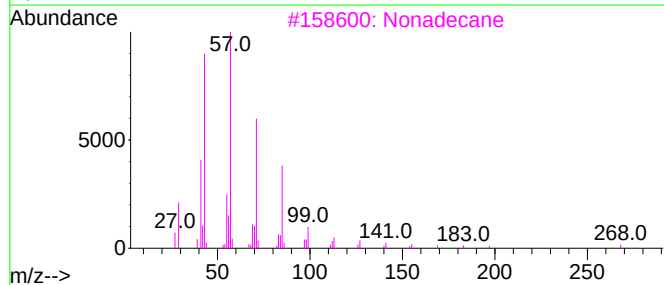
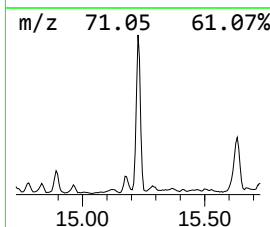
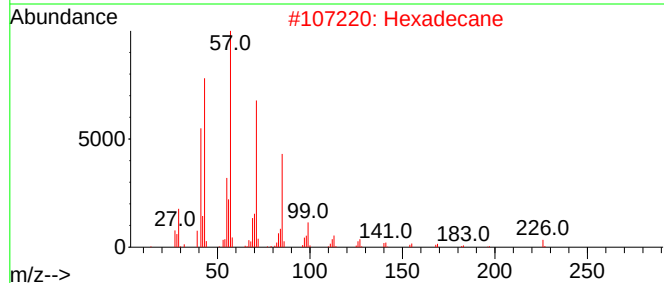
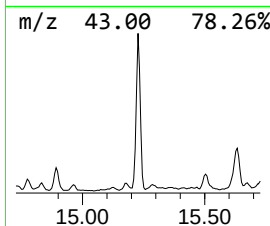
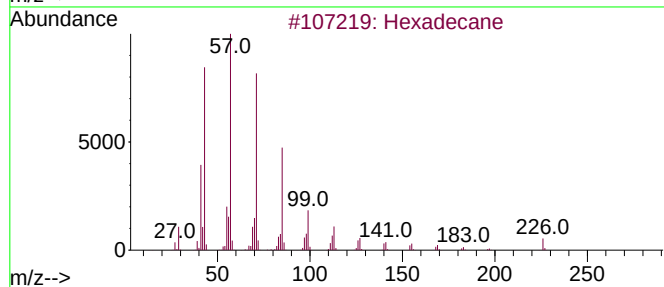
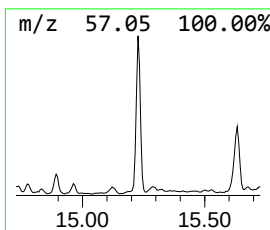
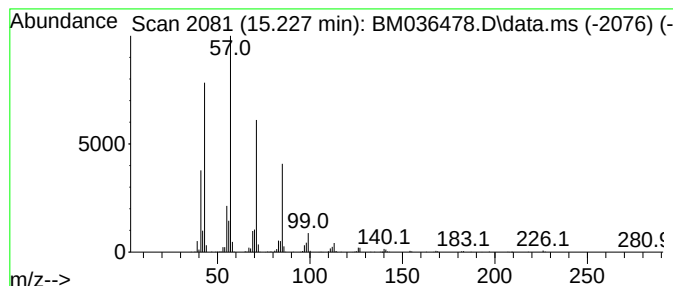
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.227	5.46 ng/ul	719139	Acenaphthene-d10	14.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	94
2			Hexadecane	226	C16H34	000544-76-3	94
3			Nonadecane	268	C19H40	000629-92-5	91
4			Heptadecane	240	C17H36	000629-78-7	90
5			Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036478.D
Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

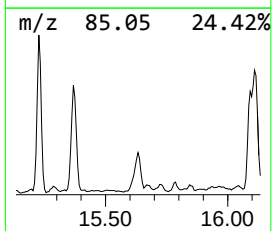
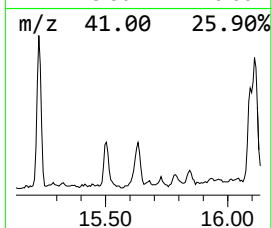
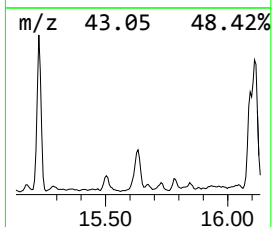
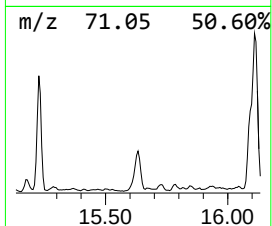
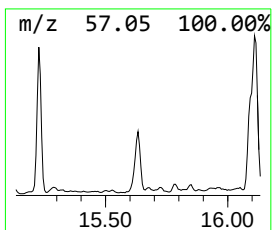
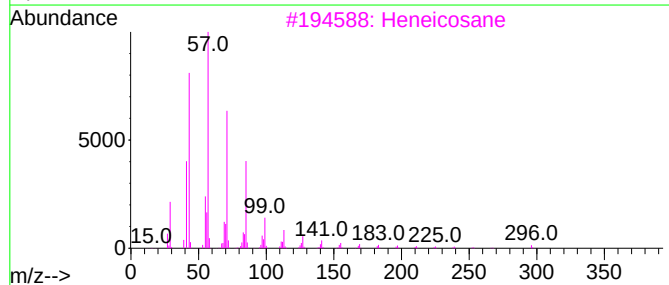
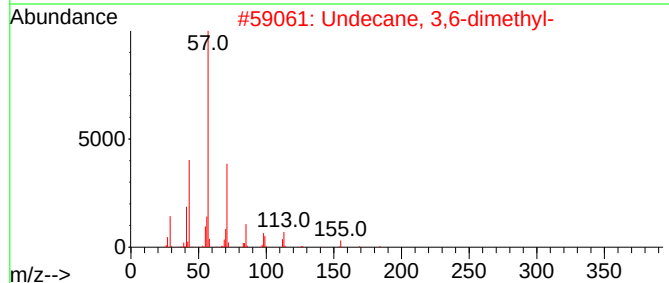
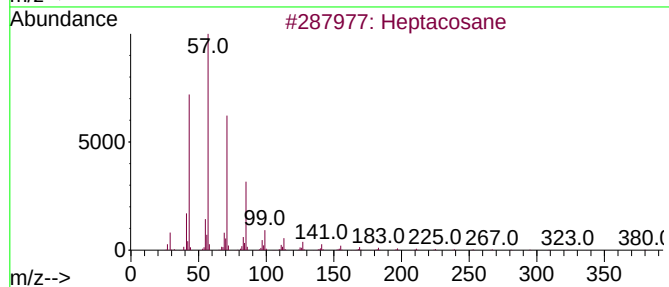
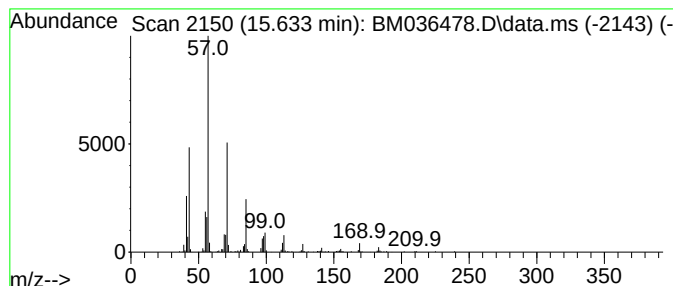
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.633	2.76 ng/ul	363902	Acenaphthene-d10		14.374	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	86
2	Undecane, 3,6-dimethyl-		184	C13H28	017301-28-9	81
3	Heneicosane		296	C21H44	000629-94-7	72
4	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	64
5	Dodecane, 3-methyl-		184	C13H28	017312-57-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036478.D
Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COKW0

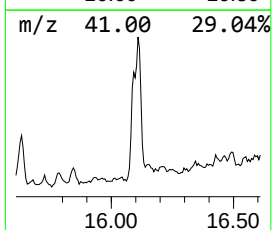
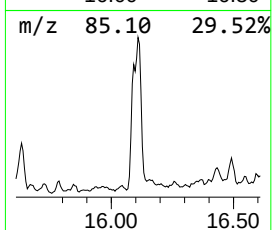
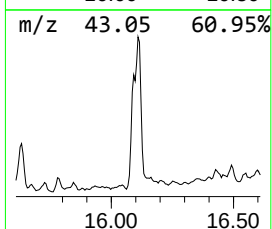
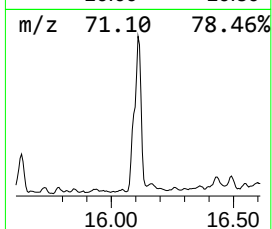
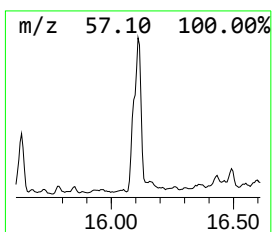
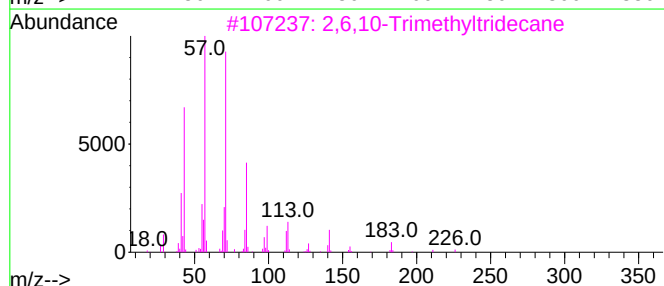
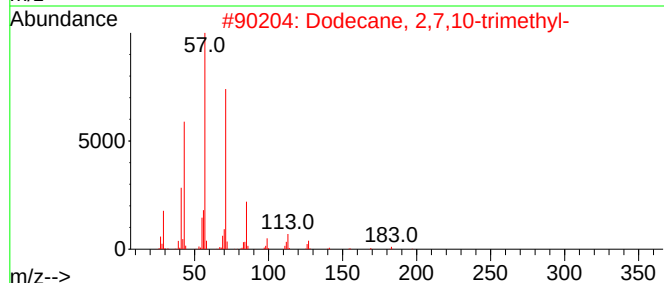
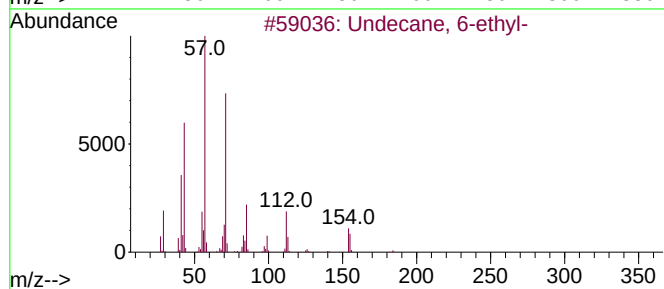
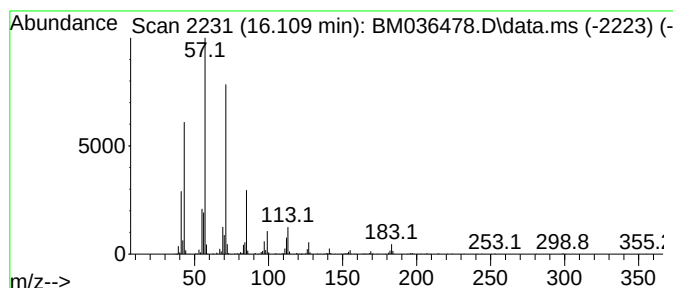
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.109	8.97 ng/ul	1454330	Phenanthrene-d10	17.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 6-ethyl-	184	C13H28	017312-60-6	87
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
3			2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	86
4			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	86
5			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036478.D
Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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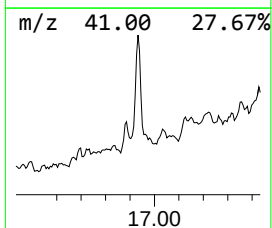
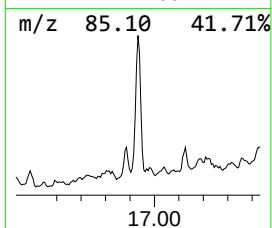
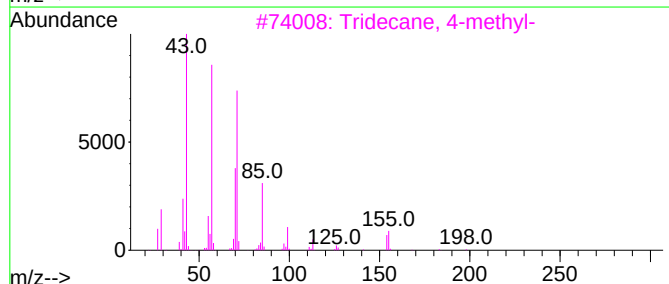
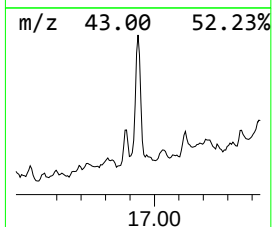
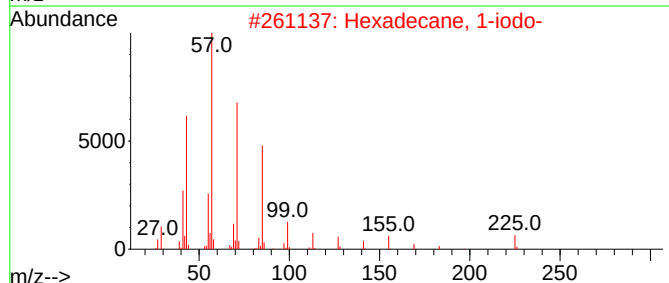
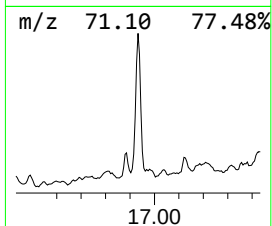
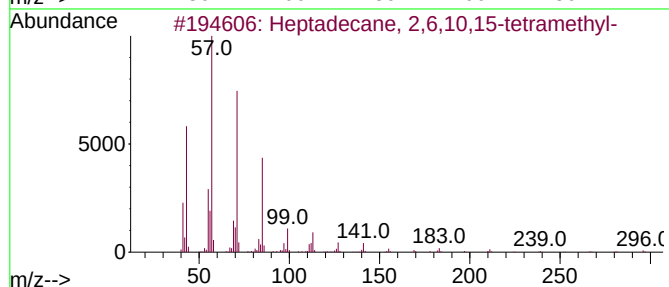
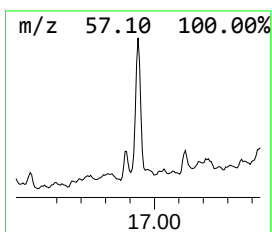
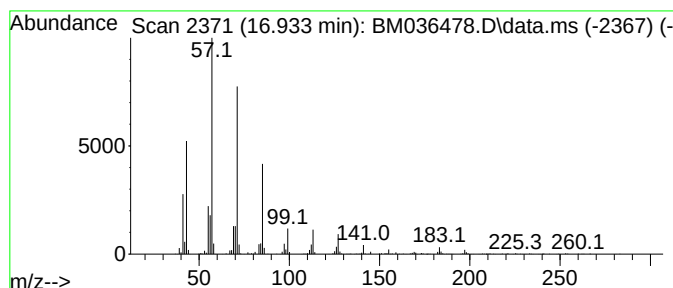
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.933	5.84 ng/ul	946994	Phenanthrene-d10	17.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	91
3			Tridecane, 4-methyl-	198	C14H30	026730-12-1	90
4			Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
5			Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
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Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

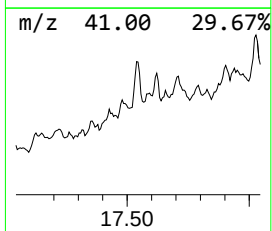
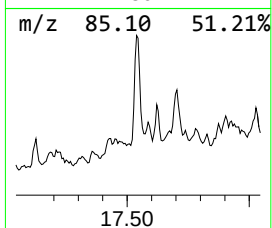
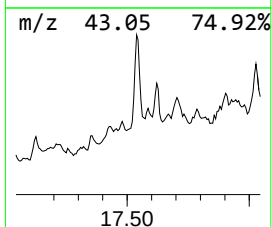
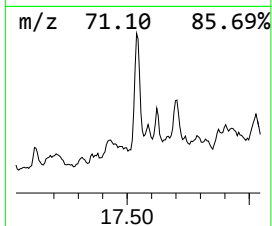
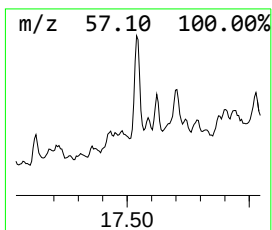
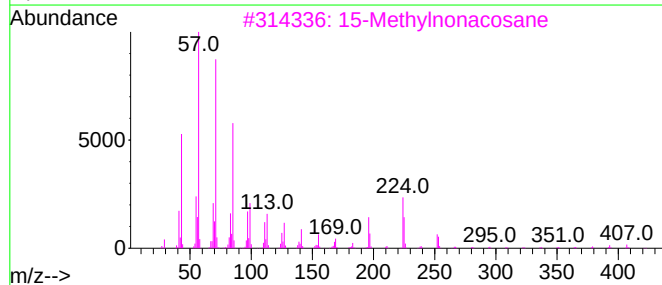
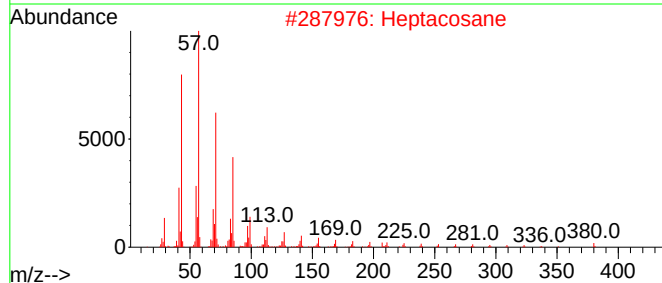
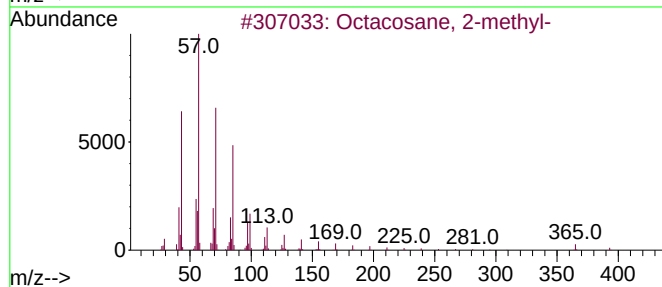
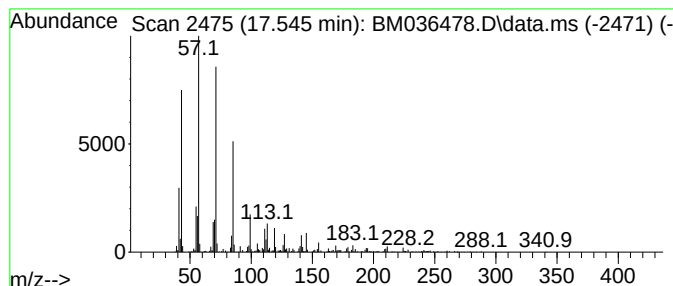
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.545	4.29 ng/ul	695623	Phenanthrene-d10		17.121	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-		408	C29H60	001560-98-1	86
2	Heptacosane		380	C27H56	000593-49-7	86
3	15-Methylnonacosane		422	C30H62	065820-60-2	86
4	Tetracosane		338	C24H50	000646-31-1	83
5	Octadecane		254	C18H38	000593-45-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036478.D
Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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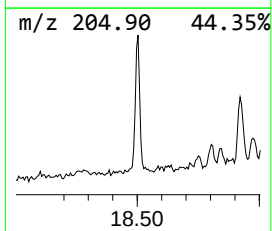
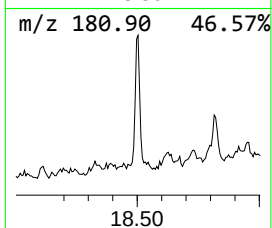
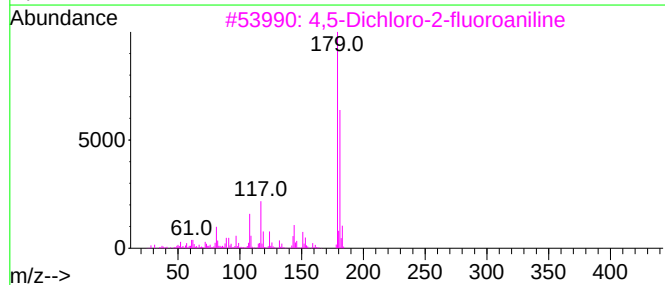
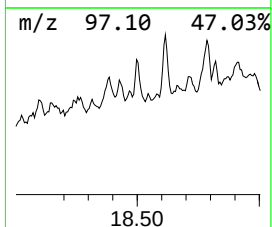
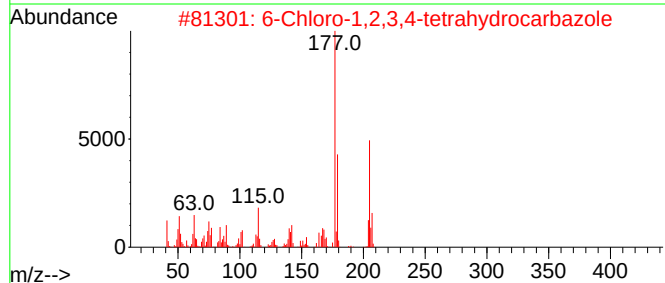
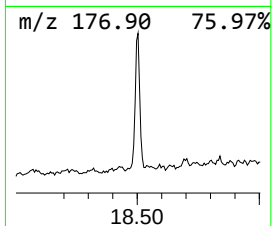
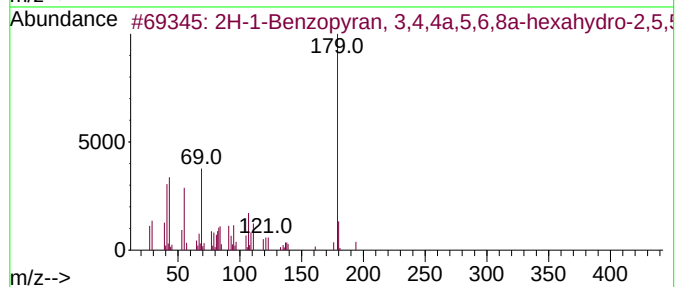
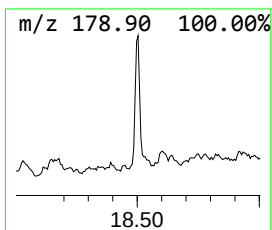
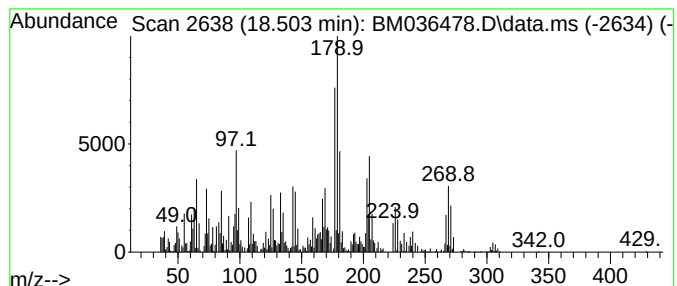
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.503	7.25 ng/ul	1174490	Phenanthrene-d10	17.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194	C13H22O	041678-32-4	41
2			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
3			4,5-Dichloro-2-fluoroaniline	179	C6H4Cl2FN	002729-36-4	20
4			1,3,5,6-Tetramethyladamantane	192	C14H24	034694-70-7	15
5			4-Butylbenzoic acid, nonyl ester	304	C20H32O2	1000292-20-6	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036478.D
Acq On : 01 Sep 2022 18:12
Operator : CG/JU
Sample : N4458-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.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
Benzene, 1-chlo...	11.139	14.7	ng/u1	1173220	2	10.527	1599630	20.0
Benzene, 1-chlo...	11.357	3.7	ng/u1	294447	2	10.527	1599630	20.0
(DEL) Alkane: S...	14.274	3.8	ng/u1	502131	3	14.374	2633360	20.0
(DEL) Alkane: S...	15.227	5.5	ng/u1	719139	3	14.374	2633360	20.0
(DEL) Alkane: S...	15.633	2.8	ng/u1	363902	3	14.374	2633360	20.0
(DEL) Alkane: S...	16.109	9.0	ng/u1	1454330	4	17.121	3241500	20.0
(DEL) Alkane: S...	16.933	5.8	ng/u1	946994	4	17.121	3241500	20.0
(DEL) Alkane: S...	17.545	4.3	ng/u1	695623	4	17.121	3241500	20.0
unknown-01	18.503	7.3	ng/u1	1174490	4	17.121	3241500	20.0