

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049706.D
 Acq On : 7 Aug 2021 15:24
 Operator : CG/JU
 Sample : M3208-07
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C00E7

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_G\METHODS\SFAM-EPA-BG080421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049706.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.078	3	7	15	rVB2	78665	169775	31.22%	2.676%
2	3.161	16	21	36	rVB	280169	543880	100.00%	8.572%
3	3.431	63	67	72	rBV6	3539	6190	1.14%	0.098%
4	4.576	261	262	266	rVB4	2560	2361	0.43%	0.037%
5	5.598	432	436	438	rBV4	1247	1650	0.30%	0.026%
6	5.845	476	478	482	rBV2	1577	2158	0.40%	0.034%
7	6.039	508	511	515	rBV5	2405	3196	0.59%	0.050%
8	7.196	702	708	716	rBV	63497	119243	21.92%	1.879%
9	7.361	730	736	744	rBV	43543	74410	13.68%	1.173%
10	7.572	765	772	782	rBV	63718	115277	21.20%	1.817%
11	7.655	782	786	789	rBV4	4586	6333	1.16%	0.100%
12	8.042	845	852	861	rVB	97693	175521	32.27%	2.766%
13	8.589	943	945	948	rVB2	1487	1542	0.28%	0.024%
14	8.753	967	973	979	rVB2	51020	94044	17.29%	1.482%
15	8.824	983	985	990	rVB3	2407	1772	0.33%	0.028%
16	8.882	990	995	999	rVB4	3338	5379	0.99%	0.085%
17	9.211	1044	1051	1058	rBV2	67563	124602	22.91%	1.964%
18	9.270	1058	1061	1065	rVB3	5110	7457	1.37%	0.118%
19	9.687	1129	1132	1135	rBV	1240	1608	0.30%	0.025%
20	9.940	1167	1175	1183	rBV2	63297	122372	22.50%	1.929%
21	10.263	1228	1230	1231	rBV	1822	1581	0.29%	0.025%
22	10.480	1261	1267	1274	rBV2	80934	150648	27.70%	2.374%
23	10.627	1286	1292	1299	rBV2	26896	48482	8.91%	0.764%
24	10.862	1322	1332	1338	rBV	137072	263859	48.51%	4.159%
25	10.997	1350	1355	1364	rBV2	41709	84328	15.50%	1.329%
26	11.620	1457	1461	1462	rBV2	1216	1695	0.31%	0.027%
27	11.678	1470	1471	1475	rBV2	1737	1959	0.36%	0.031%
28	11.767	1479	1486	1487	rBV2	1706	2669	0.49%	0.042%
29	11.878	1499	1505	1509	rBV2	8027	14301	2.63%	0.225%
30	12.266	1564	1571	1578	rBV4	9146	15610	2.87%	0.246%
31	12.389	1588	1592	1594	rBV2	1374	1811	0.33%	0.029%
32	12.448	1599	1602	1606	rVV4	2160	2927	0.54%	0.046%
33	12.489	1606	1609	1610	rVV4	2096	1861	0.34%	0.029%
34	12.519	1610	1614	1620	rVB2	10352	17555	3.23%	0.277%
35	12.736	1647	1651	1659	rVB2	5909	12388	2.28%	0.195%
36	12.812	1659	1664	1667	rBV2	1650	3139	0.58%	0.049%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049706.D
 Acq On : 7 Aug 2021 15:24
 Operator : CG/JU
 Sample : M3208-07
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C00E7

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_G\METHODS\SFAM-EPA-BG080421.M
 Title : SVOA CALIBRATION

37	13.218	1731	1733	1738	rVB2	4734	6044	1.11%	0.095%
38	13.288	1742	1745	1750	rBV2	1391	2091	0.38%	0.033%
39	13.511	1775	1783	1790	rVV4	13698	28357	5.21%	0.447%
40	13.564	1790	1792	1798	rBV3	1586	2693	0.50%	0.042%
41	13.617	1798	1801	1804	rBV2	1536	1833	0.34%	0.029%
42	13.852	1837	1841	1843	rBV	4736	7121	1.31%	0.112%
43	13.970	1856	1861	1862	rBV	2699	3031	0.56%	0.048%
44	14.011	1862	1868	1873	rBV4	9790	16284	2.99%	0.257%
45	14.093	1874	1882	1890	rVV	166730	266072	48.92%	4.194%
46	14.163	1890	1894	1897	rVB2	6178	8695	1.60%	0.137%
47	14.210	1897	1902	1905	rBV5	2737	3691	0.68%	0.058%
48	14.251	1905	1909	1912	rBV3	3255	4563	0.84%	0.072%
49	14.387	1925	1932	1941	rVB	172467	275086	50.58%	4.336%
50	14.580	1960	1965	1970	rVB2	15311	20440	3.76%	0.322%
51	14.692	1973	1984	1991	rBV	222202	352974	64.90%	5.563%
52	14.763	1991	1996	1999	rVB4	2542	3762	0.69%	0.059%
53	14.898	2013	2019	2028	rBV2	37993	74551	13.71%	1.175%
54	14.974	2031	2032	2036	rVV2	2469	2403	0.44%	0.038%
55	15.092	2047	2052	2056	rVB4	11200	17446	3.21%	0.275%
56	15.162	2060	2064	2067	rBV3	4972	6495	1.19%	0.102%
57	15.262	2080	2081	2084	rVB	3269	2005	0.37%	0.032%
58	15.309	2084	2089	2093	rBV3	7018	9531	1.75%	0.150%
59	15.356	2094	2097	2104	rVB2	7197	10122	1.86%	0.160%
60	15.420	2104	2108	2112	rVB2	3506	4850	0.89%	0.076%
61	15.485	2112	2119	2121	rBV4	6272	9434	1.73%	0.149%
62	15.526	2122	2126	2132	rVB2	18835	27151	4.99%	0.428%
63	15.591	2136	2137	2140	rBV2	1834	2117	0.39%	0.033%
64	15.685	2147	2153	2159	rBV	225927	329091	60.51%	5.187%
65	15.802	2168	2173	2179	rVB2	57545	89077	16.38%	1.404%
66	16.031	2207	2212	2218	rVB6	6733	12888	2.37%	0.203%
67	16.084	2218	2221	2222	rBV2	2638	2637	0.48%	0.042%
68	16.149	2230	2232	2235	rBV4	2558	1888	0.35%	0.030%
69	16.325	2260	2262	2265	rBV	1874	1809	0.33%	0.029%
70	16.390	2269	2273	2282	rBV3	22635	40993	7.54%	0.646%
71	16.454	2282	2284	2286	rVV2	2057	1610	0.30%	0.025%
72	16.501	2290	2292	2297	rBV2	1400	2100	0.39%	0.033%
73	16.789	2339	2341	2347	rVB3	4894	7021	1.29%	0.111%
74	16.977	2366	2373	2376	rBV4	12685	19418	3.57%	0.306%
75	17.018	2376	2380	2382	rVV5	4870	6982	1.28%	0.110%
76	17.101	2390	2394	2399	rBV5	14991	18812	3.46%	0.296%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049706.D
 Acq On : 7 Aug 2021 15:24
 Operator : CG/JU
 Sample : M3208-07
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C00E7

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_G\METHODS\SFAM-EPA-BG080421.M
 Title : SVOA CALIBRATION

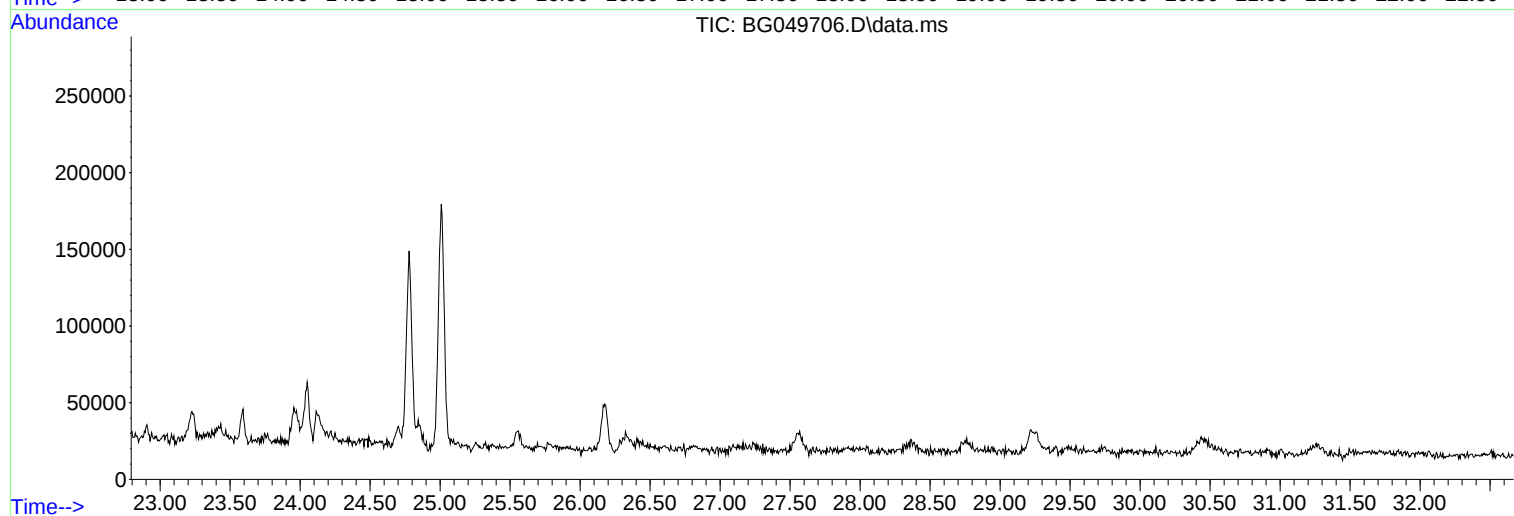
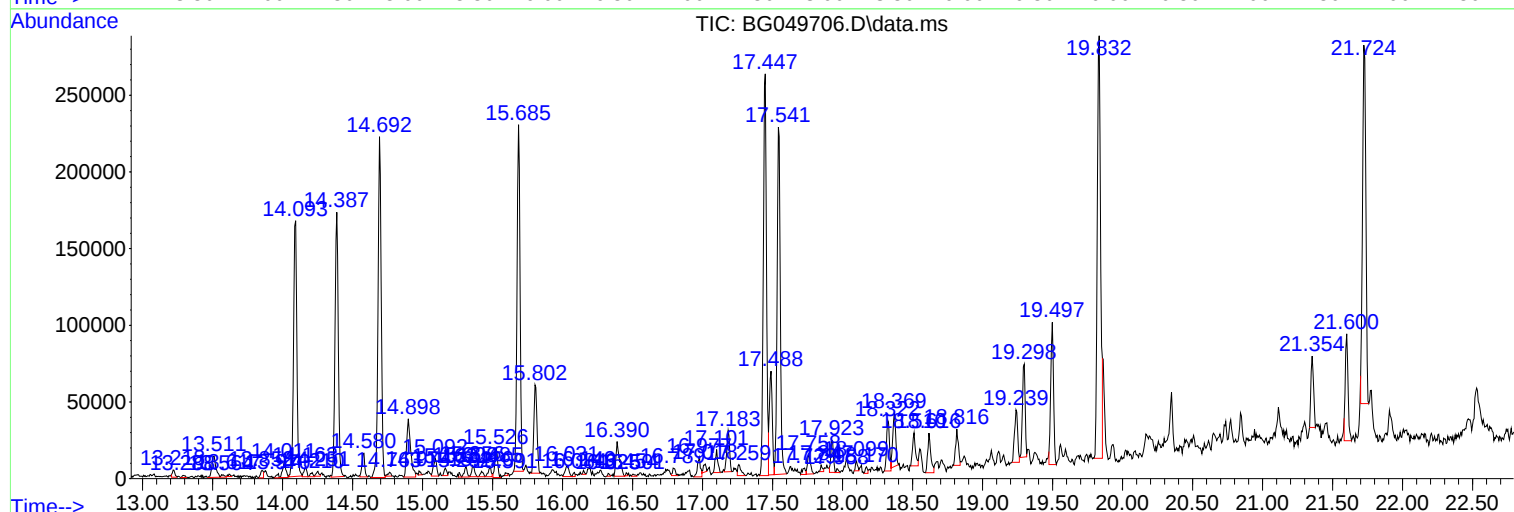
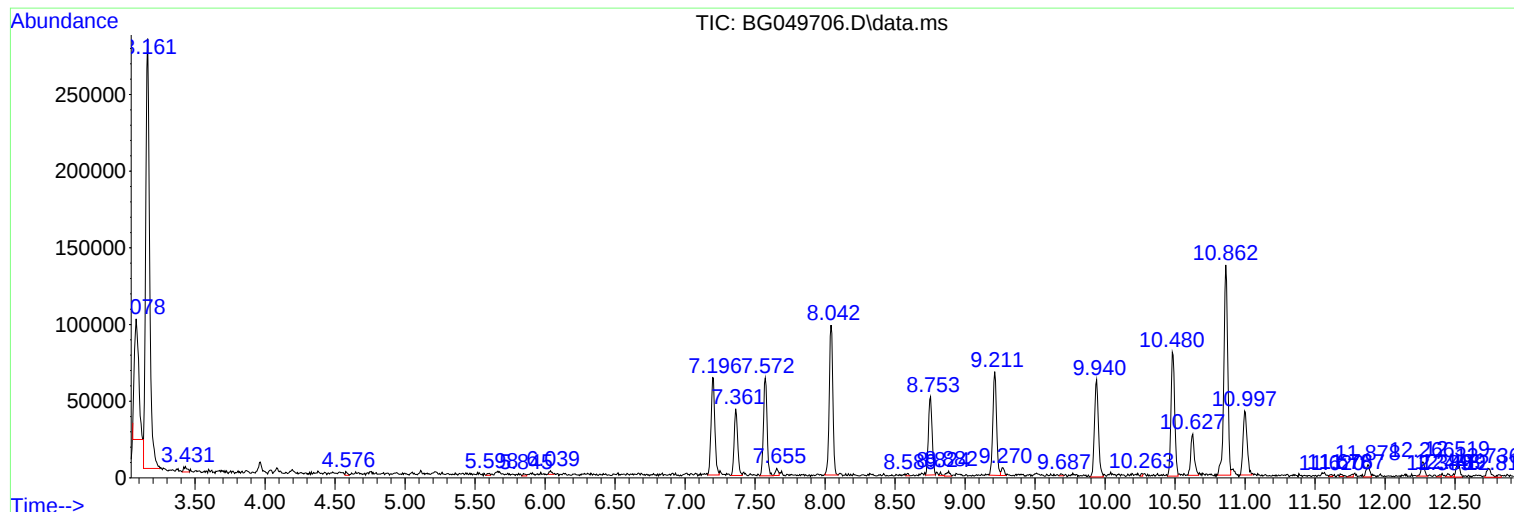
77	17.183	2403	2408	2412	rVB4	26719	37016	6.81%	0.583%
78	17.259	2419	2421	2429	rVB4	7174	11509	2.12%	0.181%
79	17.447	2446	2453	2457	rBV	262021	412522	75.85%	6.502%
80	17.488	2457	2460	2464	rVV	67654	94839	17.44%	1.495%
81	17.541	2464	2469	2479	rVB2	226260	361127	66.40%	5.692%
82	17.729	2496	2501	2503	rBV4	4151	7457	1.37%	0.118%
83	17.758	2503	2506	2512	rVB4	12614	18824	3.46%	0.297%
84	17.847	2519	2521	2524	rBV4	4267	4070	0.75%	0.064%
85	17.923	2531	2534	2538	rVB4	20985	27169	5.00%	0.428%
86	17.958	2538	2540	2544	rVB3	2514	2634	0.48%	0.042%
87	18.099	2561	2564	2567	rBV3	6245	6777	1.25%	0.107%
88	18.170	2573	2576	2578	rBV3	4276	4827	0.89%	0.076%
89	18.322	2598	2602	2606	rBV2	32778	47759	8.78%	0.753%
90	18.369	2606	2610	2615	rVB3	34641	51214	9.42%	0.807%
91	18.510	2630	2634	2639	rBV3	21599	31918	5.87%	0.503%
92	18.616	2648	2652	2658	rBV2	25887	39084	7.19%	0.616%
93	18.816	2682	2686	2690	rBV2	23714	32788	6.03%	0.517%
94	19.239	2753	2758	2762	rBV5	34241	53775	9.89%	0.848%
95	19.298	2763	2768	2771	rVV3	60498	79746	14.66%	1.257%
96	19.497	2798	2802	2807	rVB	92709	125258	23.03%	1.974%
97	19.832	2853	2859	2863	rBV	275596	458320	84.27%	7.224%
98	21.354	3115	3118	3122	rVB2	46580	60222	11.07%	0.949%
99	21.600	3157	3160	3166	rVB	69656	91489	16.82%	1.442%
100	21.724	3177	3181	3187	rVB	233727	377649	69.44%	5.952%

Sum of corrected areas: 6344744

Instrument :
BNA_G
ClientSampleId :
C00E7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

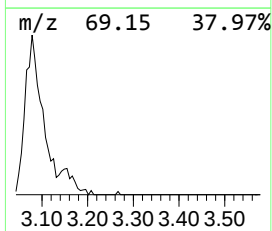
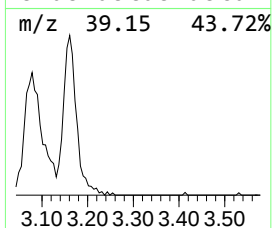
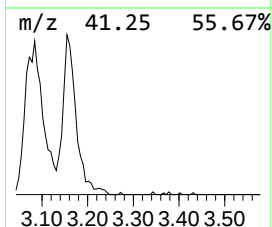
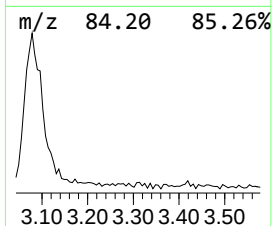
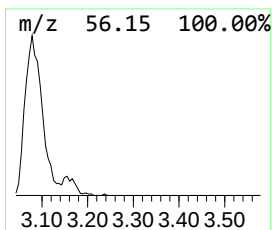
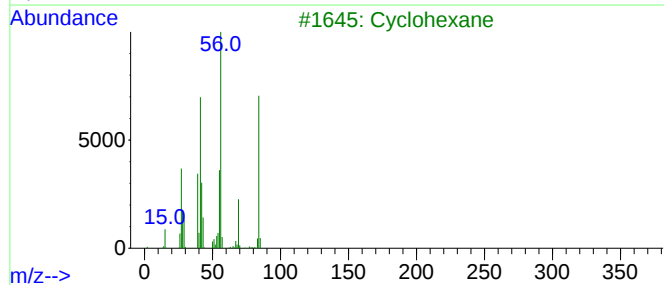
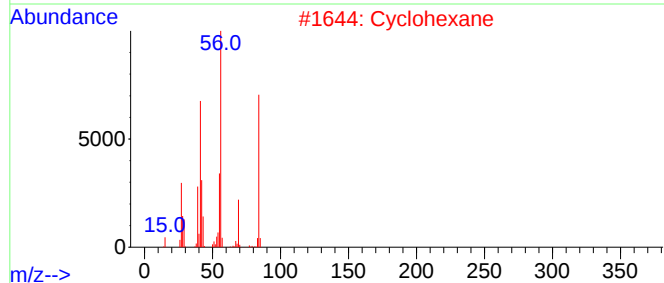
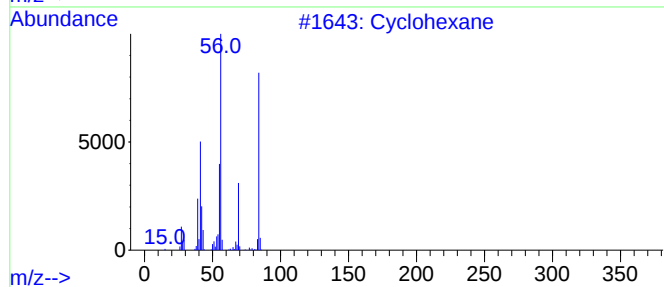
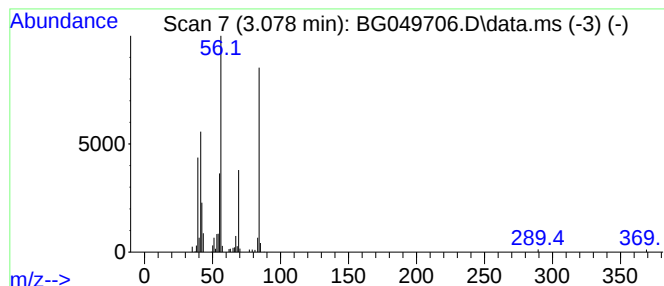
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.078 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.078	19.35 ng/ul	169775	1,4-Dichlorobenzene-d4	8.042

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	76
2		Cyclohexane	84	C6H12	000110-82-7	68
3		Cyclohexane	84	C6H12	000110-82-7	64
4		Cyclohexane	84	C6H12	000110-82-7	62
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

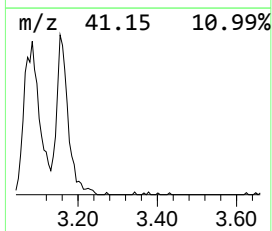
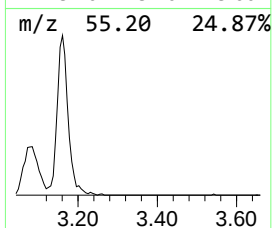
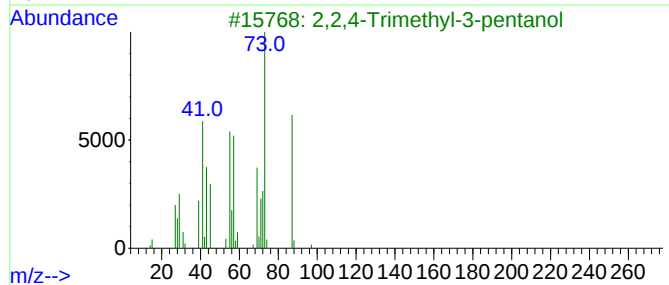
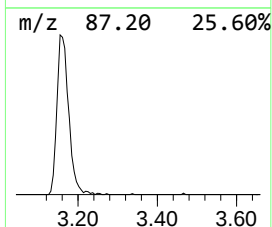
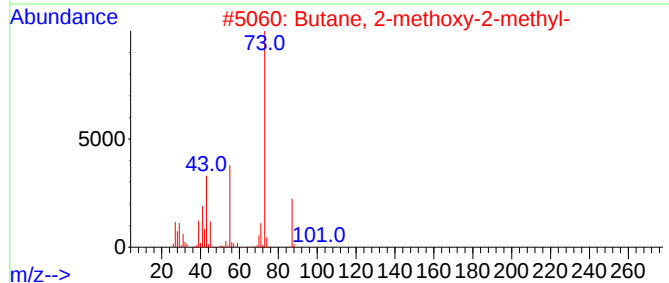
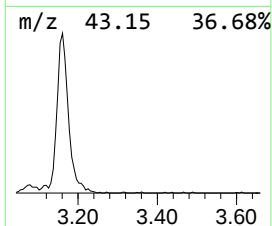
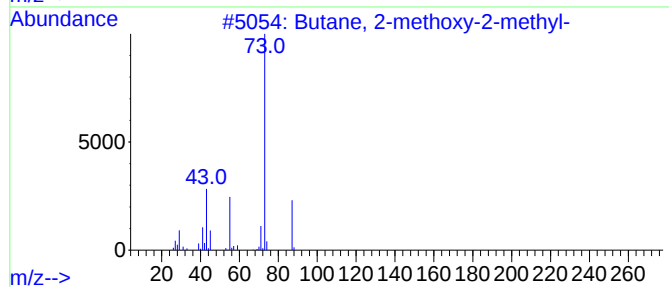
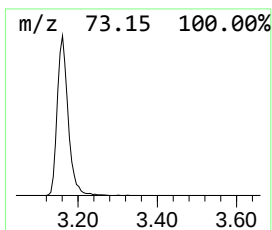
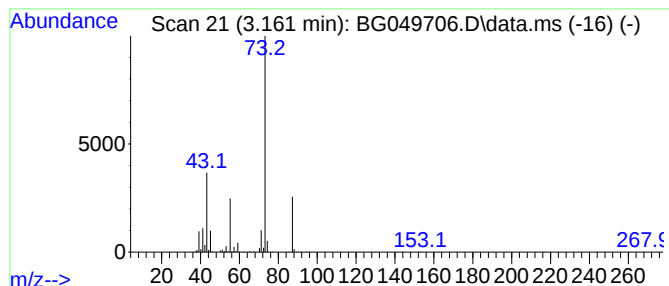
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.161	61.97 ng/ul	543880	1,4-Dichlorobenzene-d4	8.042

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

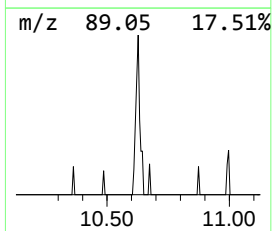
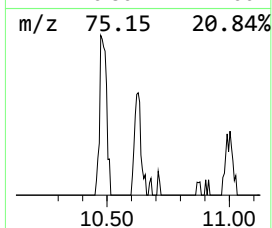
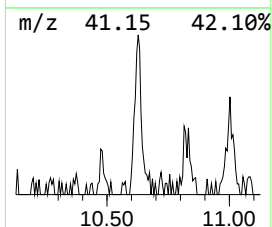
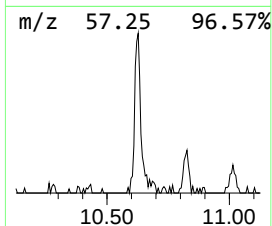
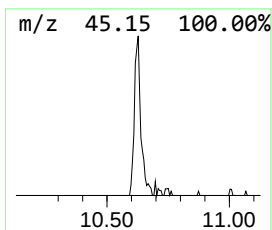
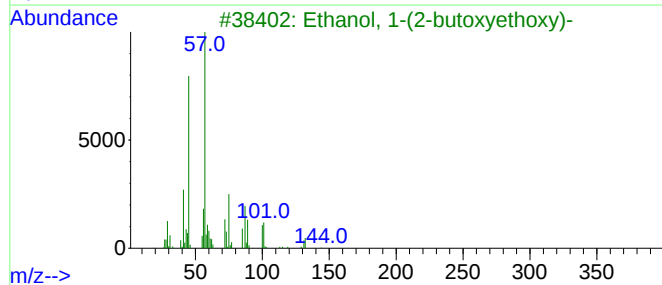
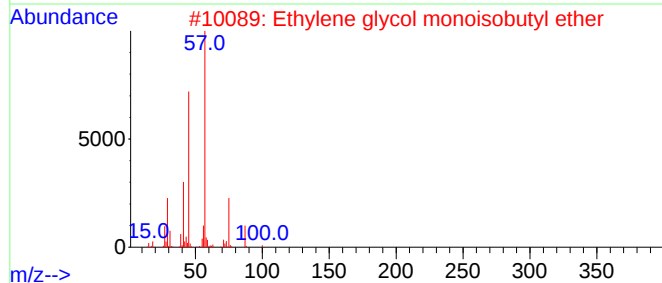
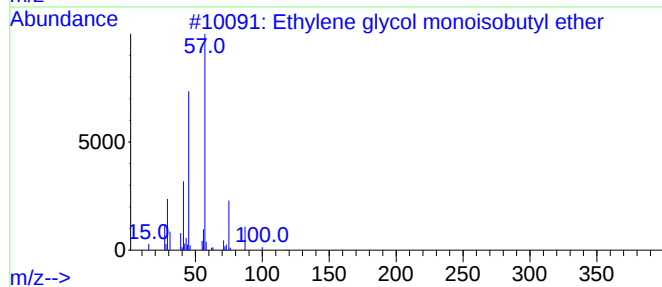
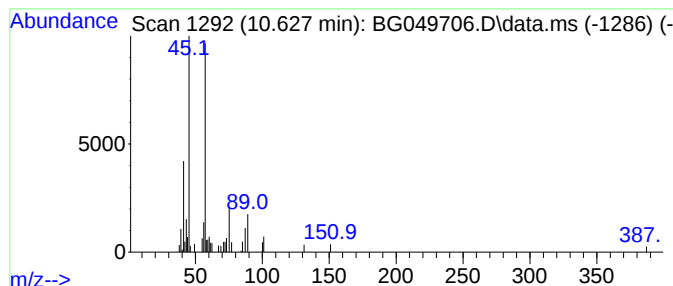
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethylene glycol monoisobuty... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.627	3.67 ng/ul	48482	Naphthalene-d8	10.862

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	50
4			Di-sec-Butyl ether	130	C8H18O	006863-58-7	50
5			Paraldehyde	132	C6H12O3	000123-63-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

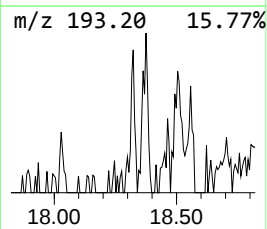
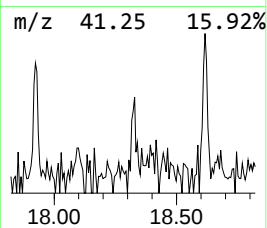
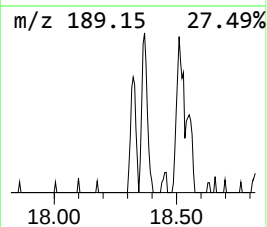
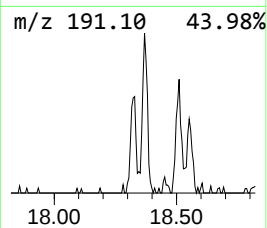
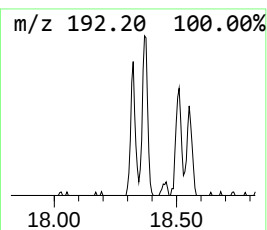
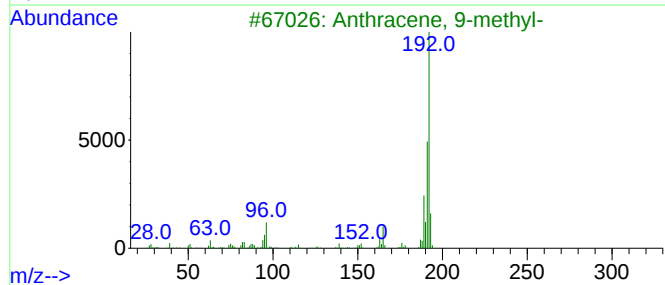
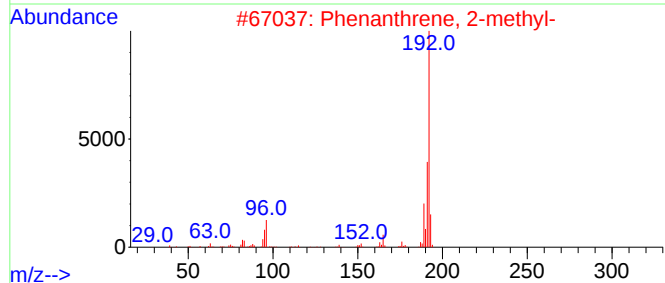
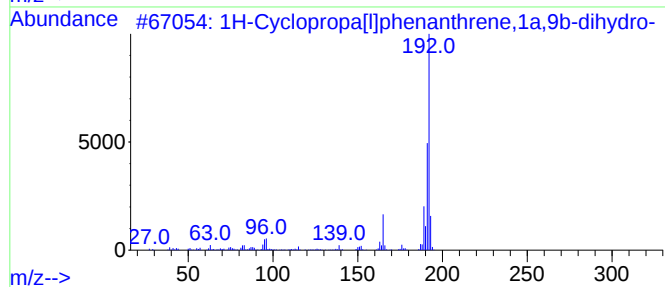
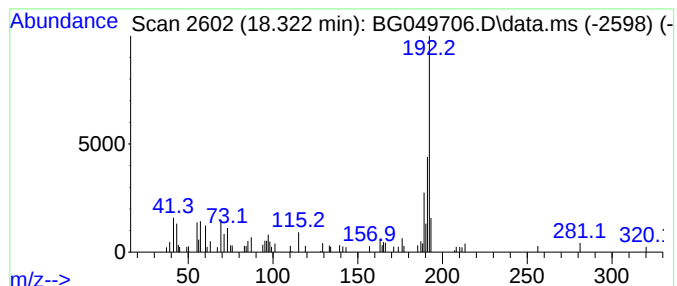
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.322	2.32 ng/ul	47759	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	76
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	76
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	76
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	76
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

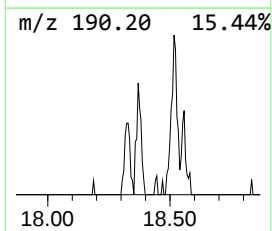
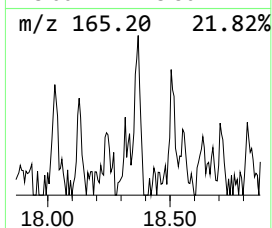
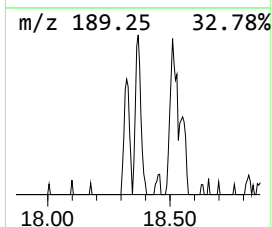
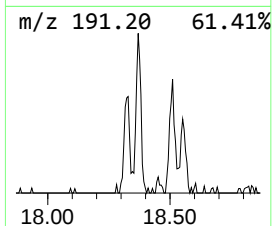
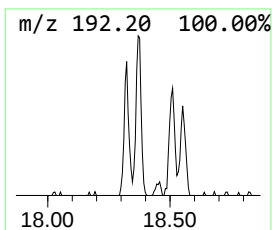
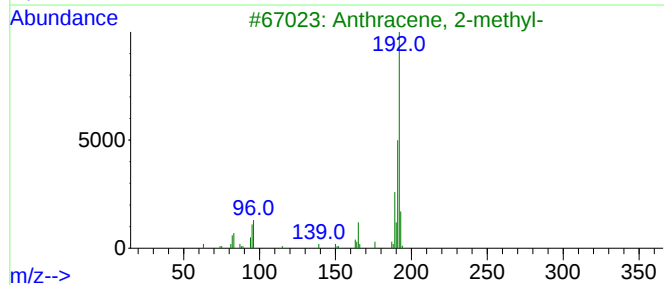
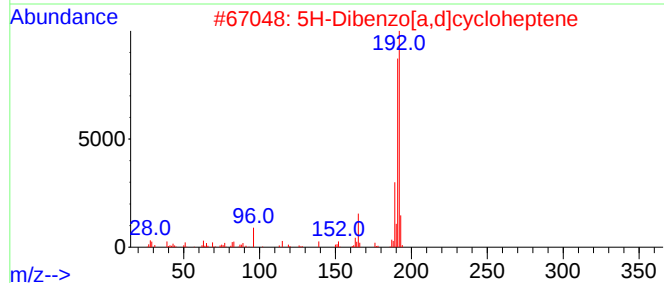
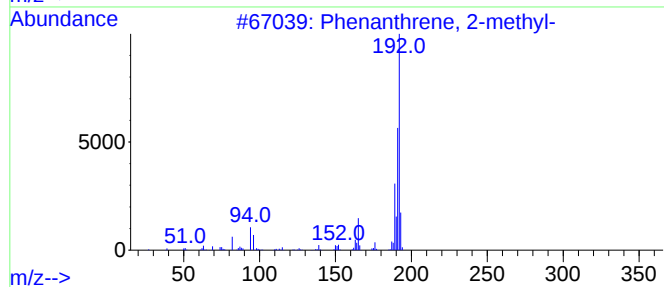
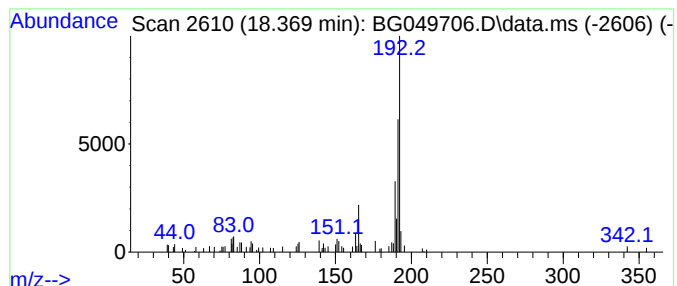
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	2.48 ng/ul	51214	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
2			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	87
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	76
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

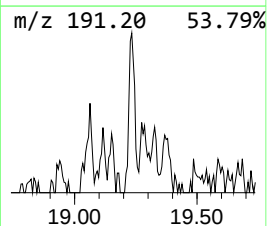
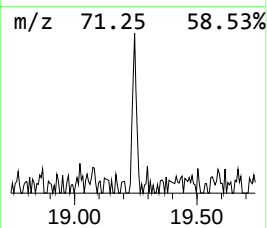
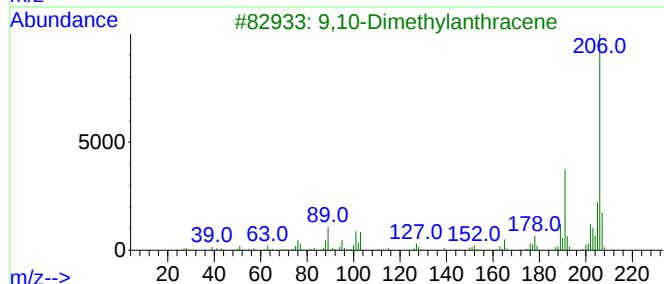
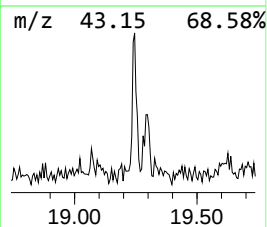
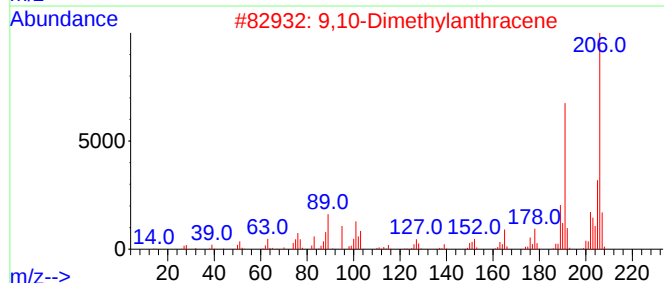
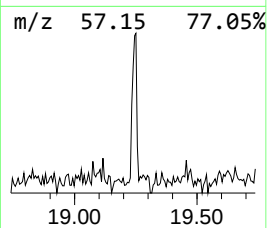
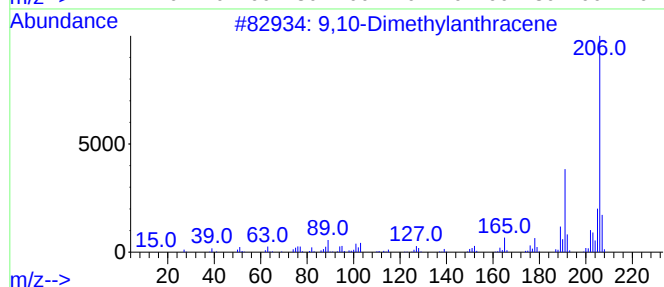
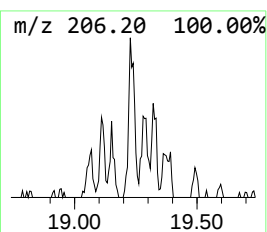
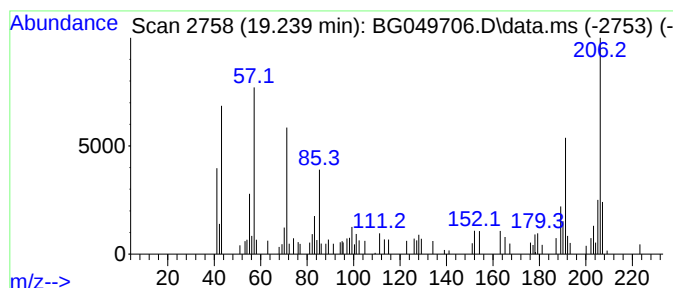
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	2.61 ng/ul	53775	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	91		
2	9,10-Dimethylantracene	206	C16H14	000781-43-1	91		
3	9,10-Dimethylantracene	206	C16H14	000781-43-1	83		
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	70		
5	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

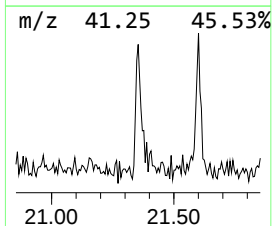
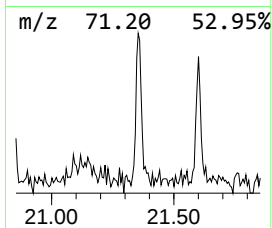
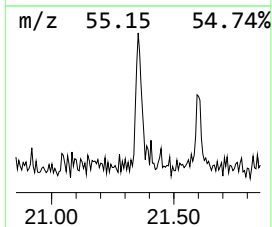
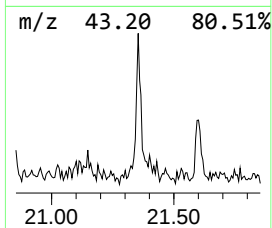
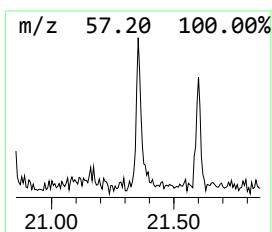
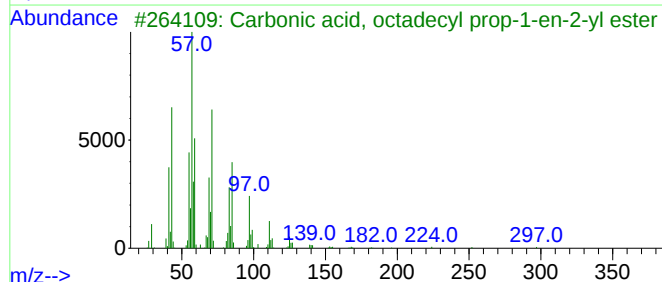
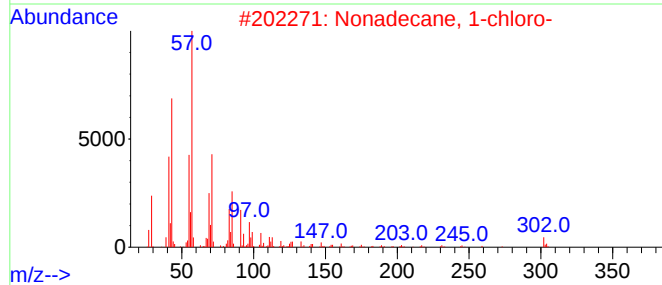
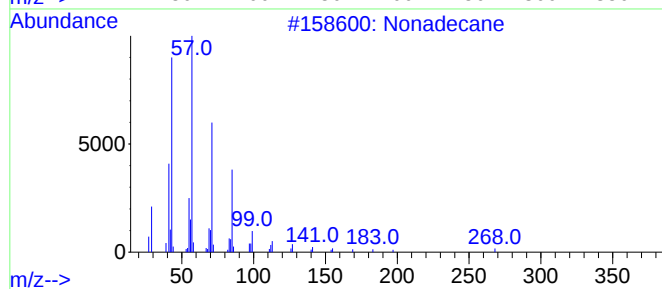
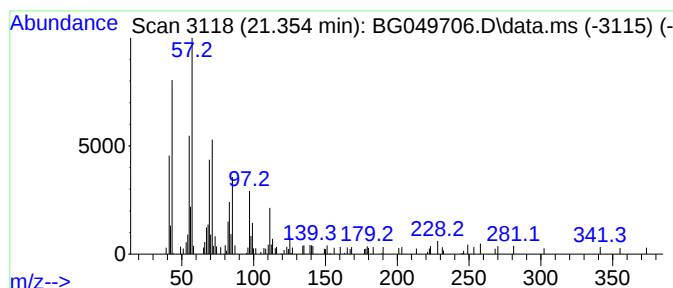
Instrument :
BNA_G
ClientSampleId :
C00E7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.354	3.19 ng/ul	60222	Chrysene-d12	21.724	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	86
2	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	76
3	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	74
4	Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	72
5	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049706.D
Acq On : 7 Aug 2021 15:24
Operator : CG/JU
Sample : M3208-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00E7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.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
(DEL) Alkane: C...	3.078	19.4	ng/ul	169775	1	8.042	175521	20.0
Butane, 2-metho...	3.161	62.0	ng/ul	543880	1	8.042	175521	20.0
Ethylene glycol...	10.627	3.7	ng/ul	48482	2	10.862	263859	20.0
1H-Cyclopropa[l...	18.322	2.3	ng/ul	47759	4	17.447	412522	20.0
Phenanthrene, 2...	18.369	2.5	ng/ul	51214	4	17.447	412522	20.0
9,10-Dimethylan...	19.239	2.6	ng/ul	53775	4	17.447	412522	20.0
Phenanthrene, 1...	19.298	3.9	ng/ul	79746	4	17.447	412522	20.0
(DEL) Alkane: S...	21.354	3.2	ng/ul	60222	5	21.724	377649	20.0