

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
 Data File : BG046786.D
 Acq On : 5 Oct 2020 16:53
 Operator : CG/JU
 Sample : L4153-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 X2N56

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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\SFAM01-EPA-BG092820.M
 Title : SVOA CALIBRATION

Signal : TIC: BG046786.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.538	29	34	39	rBV3	16245	25753	1.17%	0.066%
2	4.190	143	145	148	rVV2	3478	4031	0.18%	0.010%
3	4.225	148	151	154	rVB3	4323	4971	0.23%	0.013%
4	4.290	159	162	167	rVB4	5655	8840	0.40%	0.023%
5	4.607	214	216	221	rVB3	3270	4459	0.20%	0.011%
6	4.825	249	253	256	rVB5	5446	8496	0.39%	0.022%
7	5.206	312	318	322	rVV	12181	17443	0.80%	0.044%
8	7.098	637	640	643	rBV2	3880	4615	0.21%	0.012%
9	7.257	659	667	678	rBV2	335556	887612	40.47%	2.263%
10	7.433	689	697	704	rBV	214177	363195	16.56%	0.926%
11	7.639	724	732	739	rBV2	301973	520651	23.74%	1.328%
12	7.692	739	741	746	rVV3	4345	5789	0.26%	0.015%
13	8.109	805	812	820	rBV	216228	370222	16.88%	0.944%
14	8.802	922	930	940	rBV2	366719	629649	28.71%	1.606%
15	9.202	991	998	1004	rBV	228372	395784	18.05%	1.009%
16	9.272	1004	1010	1017	rVB	283669	505496	23.05%	1.289%
17	9.930	1116	1122	1123	rBV4	16592	23463	1.07%	0.060%
18	9.995	1124	1133	1141	rVV	271022	523862	23.89%	1.336%
19	10.065	1144	1145	1148	rVV2	3993	4873	0.22%	0.012%
20	10.101	1148	1151	1154	rVV3	8250	10729	0.49%	0.027%
21	10.136	1155	1157	1160	rVB	3640	4475	0.20%	0.011%
22	10.530	1217	1224	1233	rBV	393580	700436	31.94%	1.786%
23	10.923	1283	1291	1300	rVB	294769	515212	23.49%	1.314%
24	11.070	1304	1316	1326	rBV2	449659	1307877	59.64%	3.335%
25	11.628	1405	1411	1416	rVB4	9077	14040	0.64%	0.036%
26	11.852	1446	1449	1454	rBV4	3525	5212	0.24%	0.013%
27	12.445	1547	1550	1555	rVB4	5961	8782	0.40%	0.022%
28	12.504	1555	1560	1564	rBV2	5496	8149	0.37%	0.021%
29	13.573	1735	1742	1750	rVB	915729	1416990	64.61%	3.613%
30	13.808	1775	1782	1790	rBV2	702614	1156573	52.74%	2.949%
31	14.131	1830	1837	1842	rBV	686461	1014550	46.26%	2.587%
32	14.178	1842	1845	1851	rVB	395407	540274	24.64%	1.378%
33	14.431	1879	1888	1894	rBV	749739	1169304	53.32%	2.982%
34	14.736	1933	1940	1945	rBV2	475761	760914	34.70%	1.940%
35	14.795	1945	1950	1957	rVB	451876	705662	32.18%	1.799%
36	14.901	1961	1968	1986	rBV	359555	588963	26.86%	1.502%

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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

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM01-EPA-BG092820.M
 Title : SVOA CALIBRATION

37	15.136	2005	2008	2011	rBV4	6620	8136	0.37%	0.021%
38	15.465	2059	2064	2068	rVV5	10927	16019	0.73%	0.041%
39	15.541	2071	2077	2083	rBV	460930	622898	28.40%	1.588%
40	15.724	2100	2108	2116	rBV	1010213	1477424	67.37%	3.767%
41	15.841	2119	2128	2136	rVV3	554753	1059150	48.29%	2.701%
42	15.964	2143	2149	2152	rVB3	10764	19299	0.88%	0.049%
43	16.505	2236	2241	2246	rBV5	4883	7926	0.36%	0.020%
44	16.652	2262	2266	2269	rBV2	13999	20979	0.96%	0.053%
45	16.699	2269	2274	2279	rVB2	106424	163343	7.45%	0.417%
46	16.781	2281	2288	2295	rVB2	753673	1066507	48.63%	2.720%
47	17.016	2324	2328	2330	rVB3	4608	5393	0.25%	0.014%
48	17.122	2339	2346	2352	rBV	985968	1332509	60.76%	3.398%
49	17.263	2368	2370	2374	rVB5	3345	4330	0.20%	0.011%
50	17.474	2399	2406	2412	rBV	660773	968662	44.17%	2.470%
51	17.521	2412	2414	2416	rVV2	4374	4315	0.20%	0.011%
52	17.574	2416	2423	2427	rVV2	1073207	1725380	78.67%	4.400%
53	17.610	2427	2429	2435	rVV	397120	492510	22.46%	1.256%
54	18.033	2491	2501	2505	rBV2	44284	74208	3.38%	0.189%
55	18.432	2564	2569	2577	rBV	759541	946305	43.15%	2.413%
56	18.879	2641	2645	2650	rVB4	22558	33125	1.51%	0.084%
57	19.084	2677	2680	2682	rBV3	3970	4198	0.19%	0.011%
58	19.108	2682	2684	2690	rVB7	2948	3915	0.18%	0.010%
59	19.490	2745	2749	2754	rVV2	19126	23014	1.05%	0.059%
60	19.637	2767	2774	2775	rBV6	2884	5060	0.23%	0.013%
61	19.742	2788	2792	2796	rBV2	14225	20607	0.94%	0.053%
62	19.854	2804	2811	2821	rVV	1527204	2193087	100.00%	5.592%
63	19.930	2821	2824	2826	rVB4	4691	4127	0.19%	0.011%
64	19.954	2826	2828	2831	rBV3	4633	4649	0.21%	0.012%
65	20.136	2856	2859	2860	rBV3	5196	5476	0.25%	0.014%
66	20.248	2874	2878	2880	rBV3	4183	5934	0.27%	0.015%
67	20.389	2900	2902	2905	rBV4	4882	5572	0.25%	0.014%
68	20.506	2920	2922	2925	rBV4	4710	5212	0.24%	0.013%
69	20.547	2928	2929	2934	rBV5	4097	5013	0.23%	0.013%
70	20.765	2961	2966	2971	rBV	1000806	1186797	54.12%	3.026%
71	20.894	2983	2988	2993	rBV	134148	204812	9.34%	0.522%
72	21.640	3109	3115	3122	rVB	1196722	1577992	71.95%	4.024%
73	21.752	3127	3134	3138	rVV	723231	1210048	55.18%	3.086%
74	21.799	3138	3142	3149	rVB	623075	999691	45.58%	2.549%
75	21.957	3167	3169	3177	rVB8	18589	33366	1.52%	0.085%
76	22.228	3212	3215	3219	rBV6	11090	16567	0.76%	0.042%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 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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77	22.369	3235	3239	3243	rBV5	13386	22966	1.05%	0.059%
78	22.545	3264	3269	3272	rBV5	32937	56483	2.58%	0.144%
79	22.768	3304	3307	3311	rBV6	7722	11202	0.51%	0.029%
80	22.827	3313	3317	3319	rVV2	37476	55240	2.52%	0.141%
81	22.874	3319	3325	3332	rVB	592424	1030951	47.01%	2.629%
82	23.003	3344	3347	3354	rVB3	31652	52358	2.39%	0.134%
83	23.121	3361	3367	3369	rBV8	15302	25294	1.15%	0.064%
84	23.291	3391	3396	3402	rVB3	37871	73361	3.35%	0.187%
85	23.538	3433	3438	3443	rBV7	19245	40966	1.87%	0.104%
86	23.873	3489	3495	3496	rBV5	16302	22794	1.04%	0.058%
87	23.943	3499	3507	3518	rVB	449452	1073284	48.94%	2.737%
88	24.114	3531	3536	3546	rVB3	60923	137335	6.26%	0.350%
89	24.184	3546	3548	3550	rBV3	7829	7356	0.34%	0.019%
90	24.801	3643	3653	3660	rBV	743094	2116698	96.52%	5.397%
91	24.866	3660	3664	3675	rVB	277343	697125	31.79%	1.778%
92	25.024	3681	3691	3698	rBV	419620	1135198	51.76%	2.895%
93	25.089	3699	3702	3712	rVB3	77996	175302	7.99%	0.447%
94	26.258	3893	3901	3909	rVB4	68695	205298	9.36%	0.523%
95	26.781	3988	3990	3991	rBV2	7395	5688	0.26%	0.015%
96	26.946	4016	4018	4021	rBV4	7407	10089	0.46%	0.026%
97	27.651	4130	4138	4150	rVB4	66343	253028	11.54%	0.645%
98	28.356	4256	4258	4259	rBV2	8380	4765	0.22%	0.012%
99	29.313	4417	4421	4422	rBV4	22041	28823	1.31%	0.073%
100	29.343	4422	4426	4443	rVB6	46950	174235	7.94%	0.444%

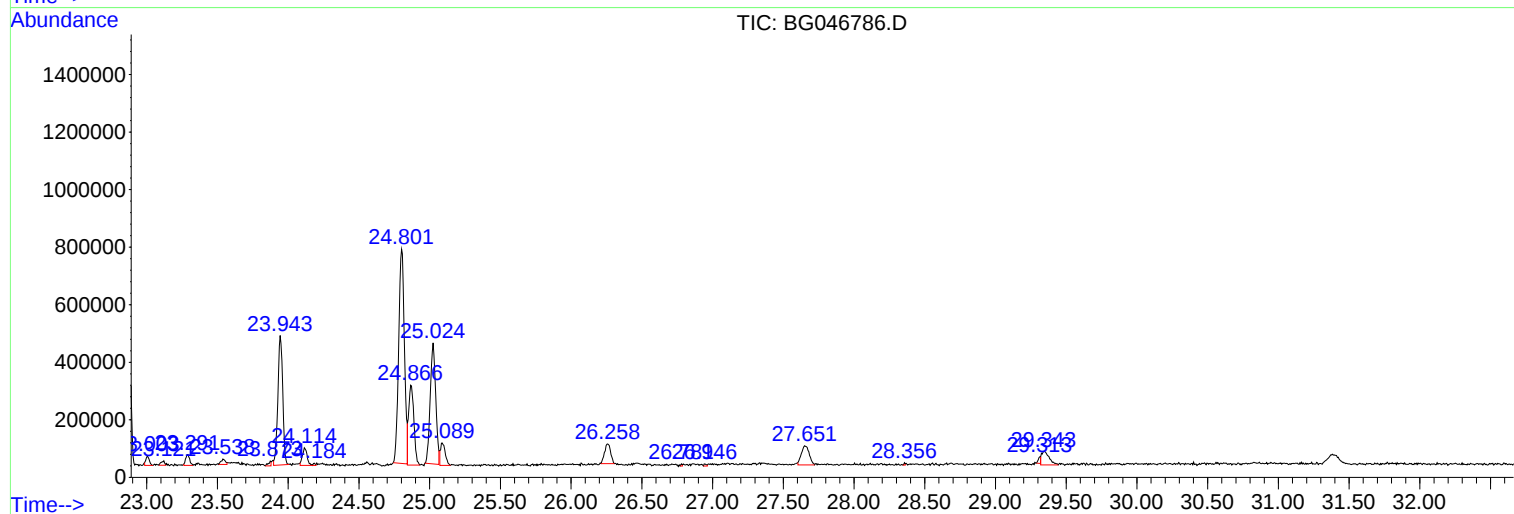
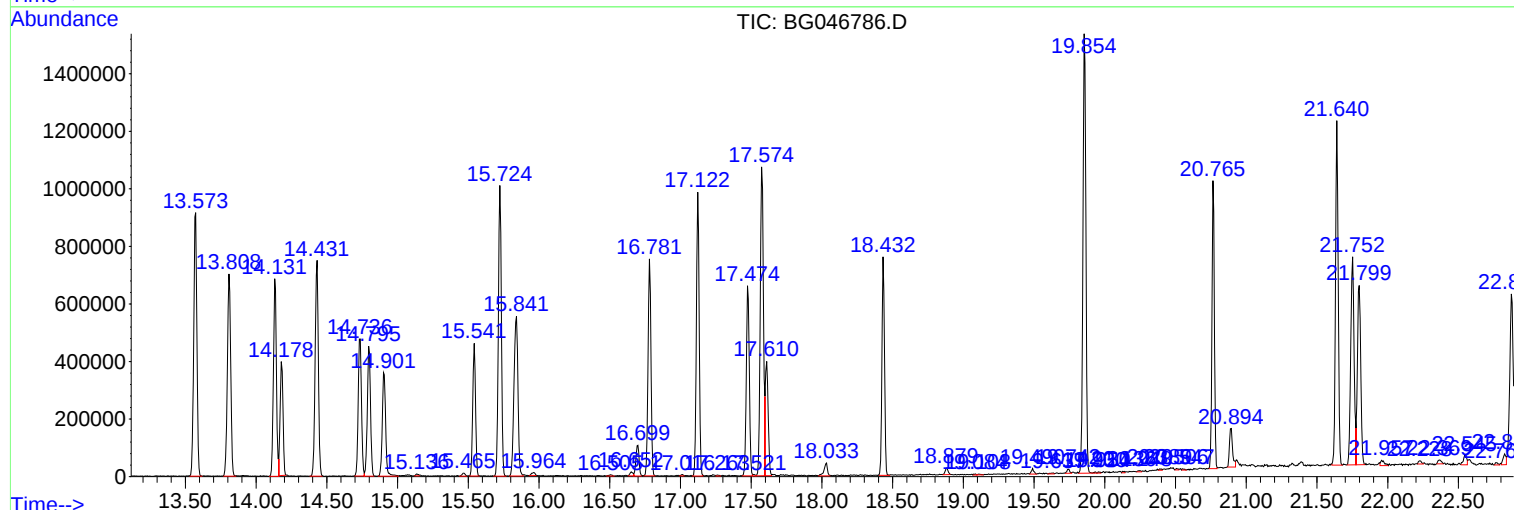
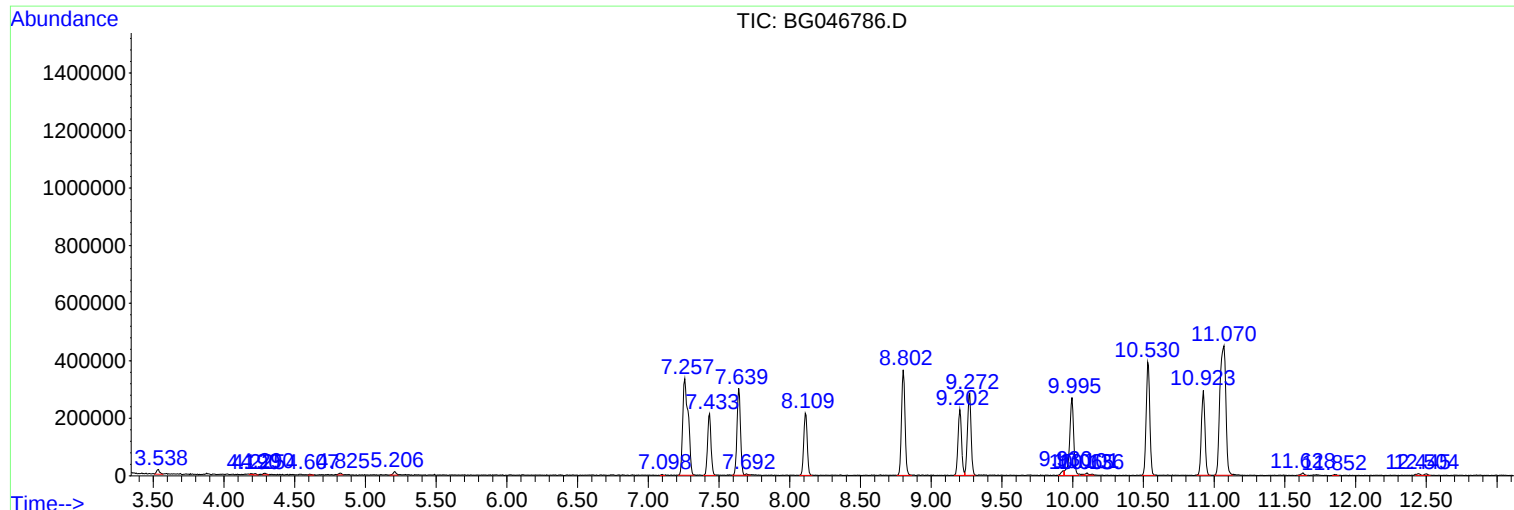
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Instrument :
BNA_G
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM01-EPA-BG092820.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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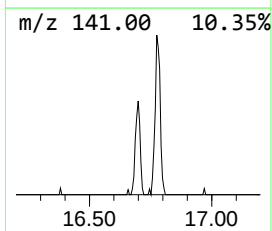
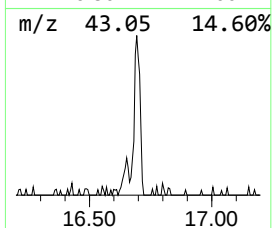
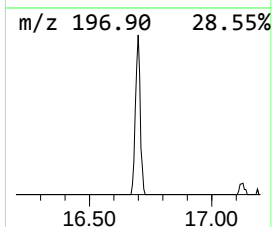
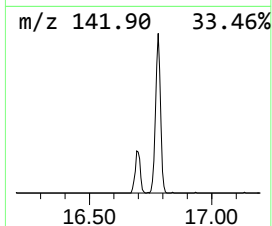
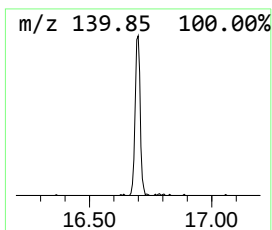
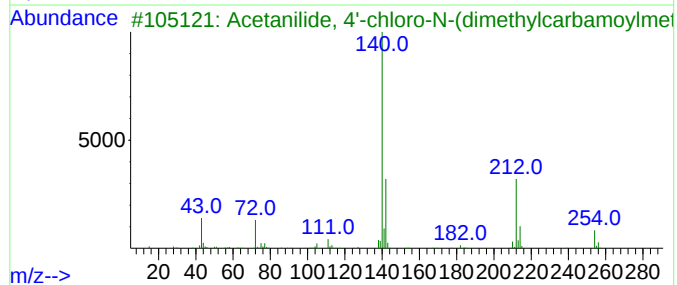
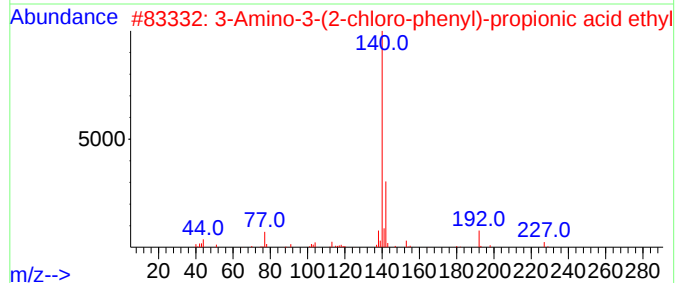
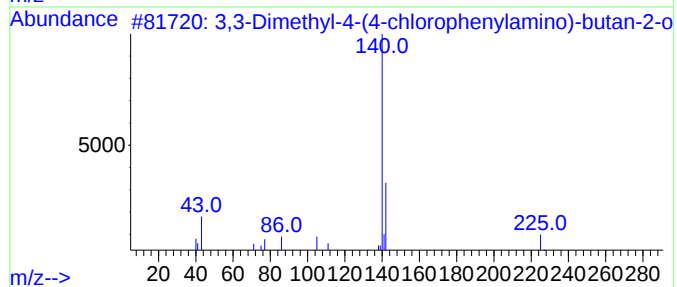
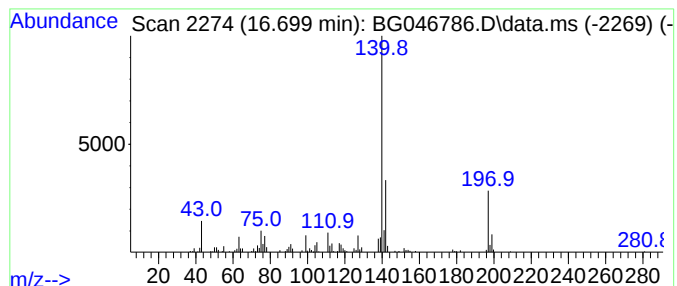
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM01-EPA-BG092820.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3,3-Dimethyl-4-(4-chlorophe... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.700	3.37 ng/ul	163343	Phenanthrene-d10	17.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,3-Dimethyl-4-(4-chlorophenylam...	225	C12H16ClNO	085937-26-4	64
2			3-Amino-3-(2-chloro-phenyl)-prop...	227	C11H14ClNO2	021464-58-4	53
3			Acetanilide, 4'-chloro-N-(dimeth...	254	C12H15ClN2O2	105900-85-4	50
4			1H-Pyrazole-4-sulfonamide, N-[(2...	299	C12H14ClN3O2S	1000320-19-0	45
5			p-Chloro-N-ethylaniline	155	C8H10ClN	013519-75-0	45



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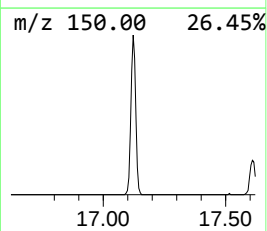
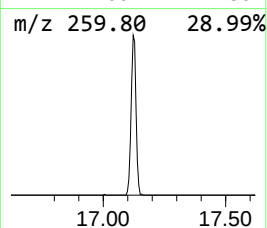
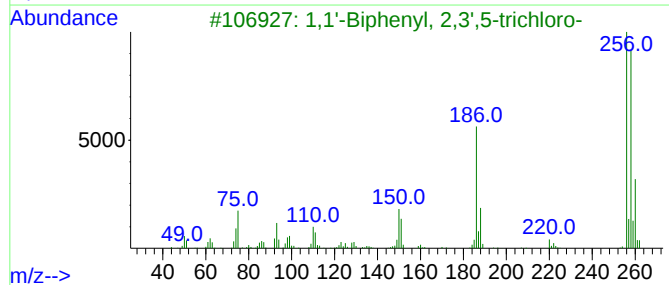
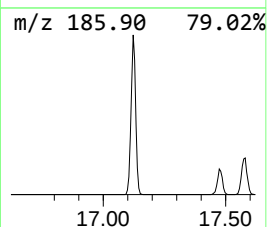
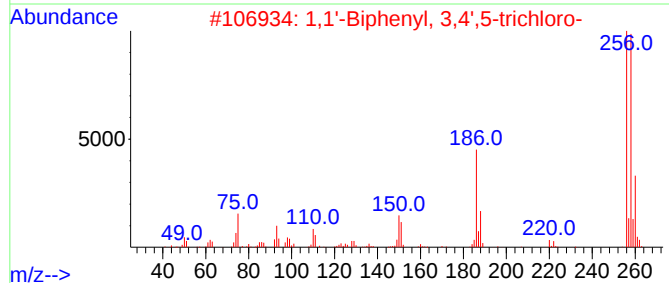
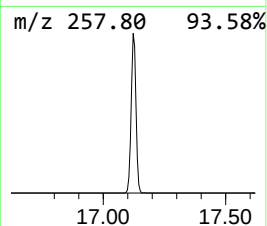
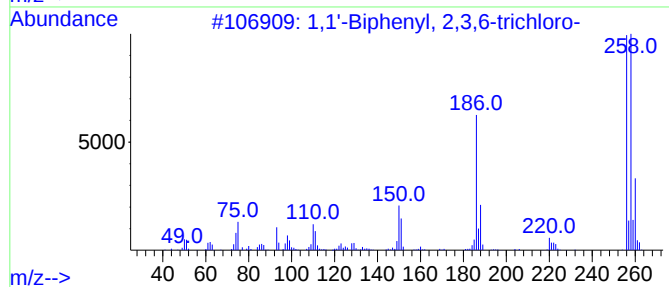
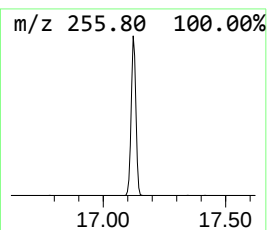
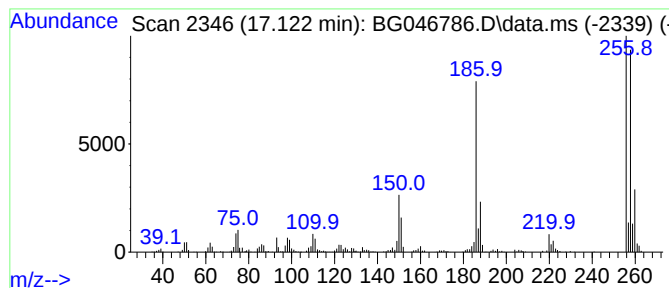
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM01-EPA-BG092820.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.120	27.51 ng/ul	1332510	Phenanthrene-d10	17.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99		
2	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99		
3	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		
4	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99		
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		



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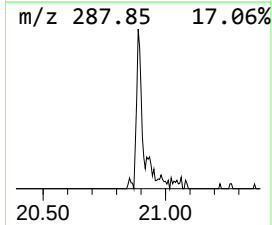
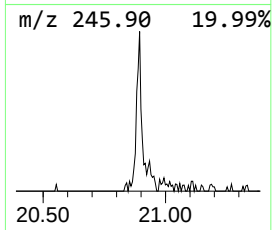
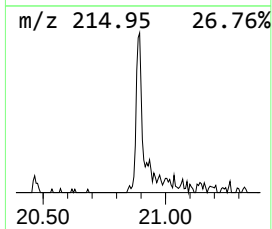
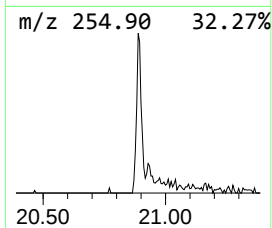
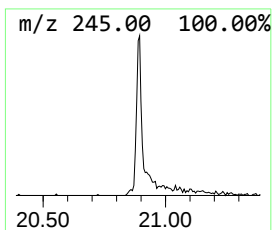
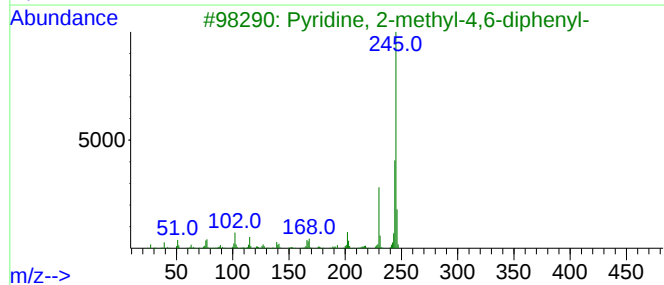
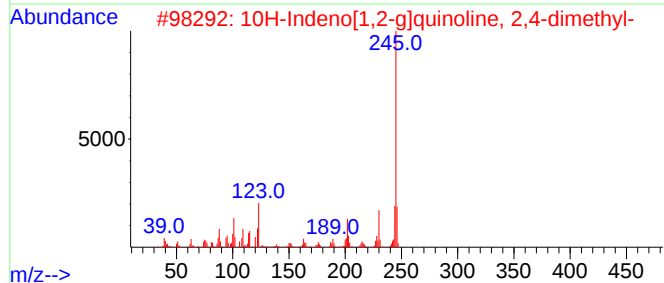
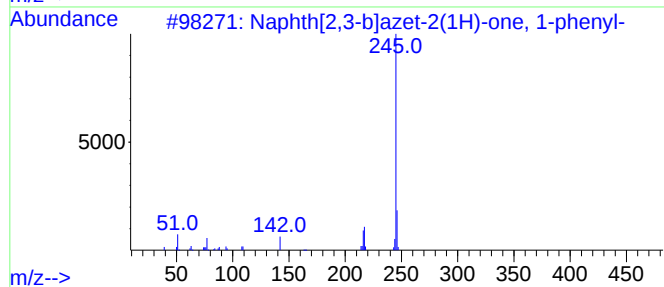
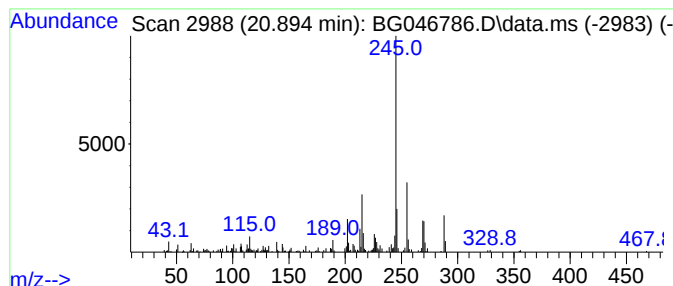
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphth[2,3-b]azet-2(1H)-one... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.890	3.39 ng/ul	204812	Chrysene-d12	21.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphth[2,3-b]azet-2(1H)-one, 1-p...	245	C17H11NO	019275-01-5	53
2			10H-Indeno[1,2-g]quinoline, 2,4-...	245	C18H15N	050519-37-4	46
3			Pyridine, 2-methyl-4,6-diphenyl-	245	C18H15N	001912-16-9	46
4			Benzenamine, N,N-diphenyl-	245	C18H15N	000603-34-9	46
5			Alpha-(4-cinnolinyl)-alpha-pheny...	245	C16H11N3	018592-05-7	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
Operator : CG/JU
Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

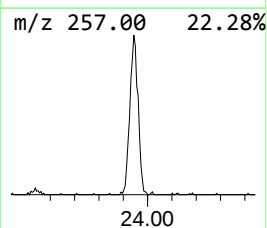
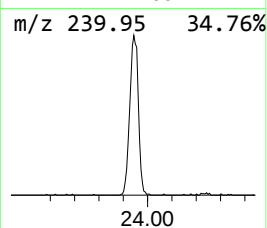
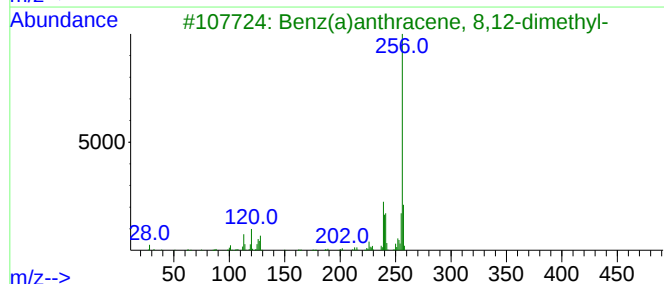
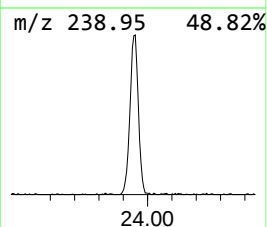
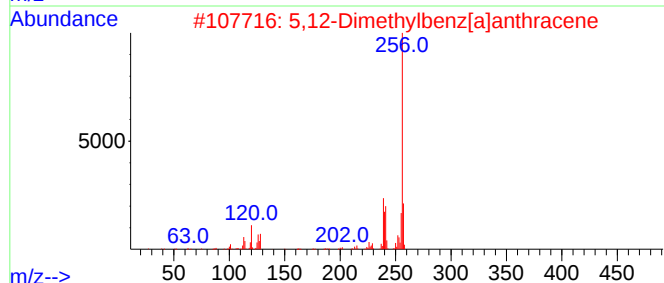
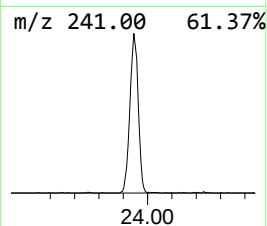
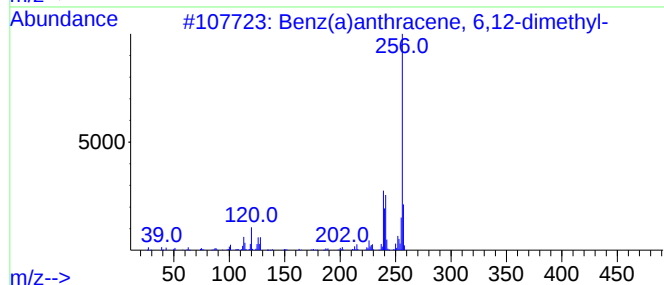
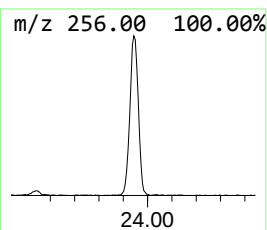
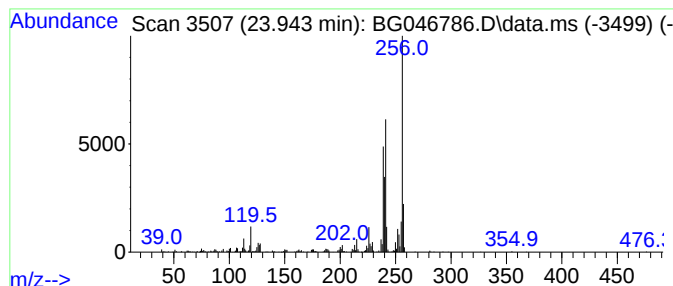
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benz(a)anthracene, 6,12-dim... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.940	18.91 ng/ul	1073280	Perylene-d12	25.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz(a)anthracene, 6,12-dimethyl-	256	C20H16	000568-81-0	98
2	5,12-Dimethylbenz[a]anthracene	256	C20H16	021297-22-3	96
3	Benz(a)anthracene, 8,12-dimethyl-	256	C20H16	020627-31-0	96
4	Benz(a)anthracene, 6,7-dimethyl-	256	C20H16	020627-28-5	96
5	Benz[a]anthracene, 1,12-dimethyl-	256	C20H16	000313-74-6	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
Operator : CG/JU
Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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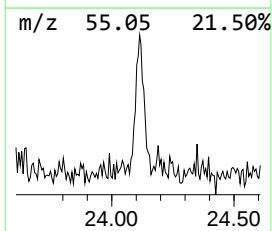
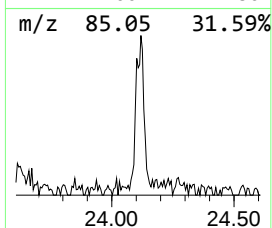
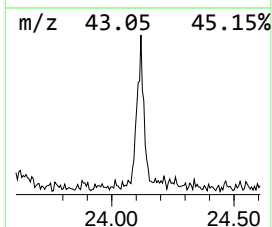
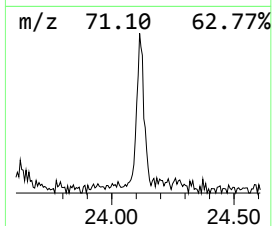
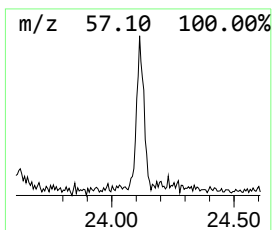
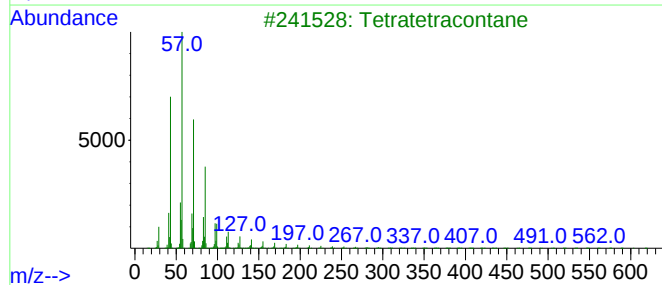
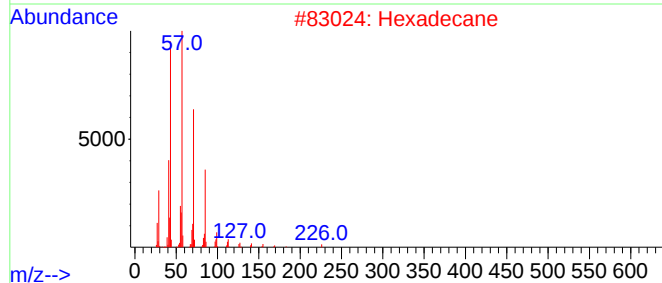
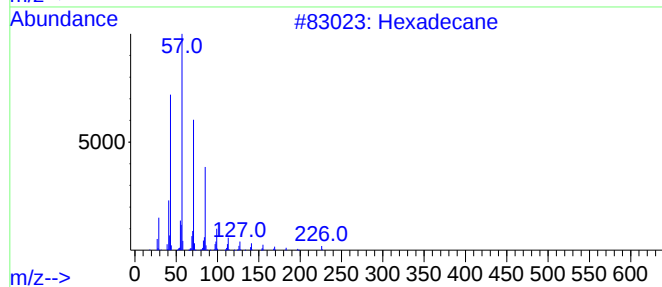
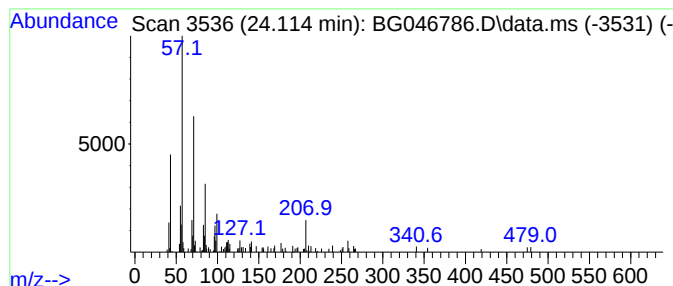
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.110	2.42 ng/ul	137335	Perylene-d12	25.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	95
2			Hexadecane	226	C16H34	000544-76-3	70
3			Tetratetracontane	619	C44H90	007098-22-8	68
4			Eicosane	282	C20H42	000112-95-8	64
5			5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
Operator : CG/JU
Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

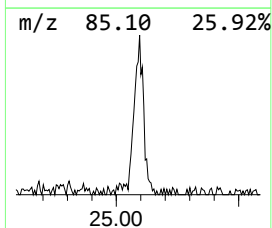
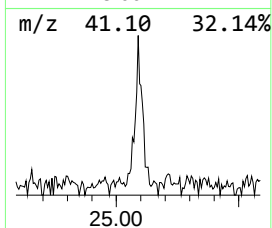
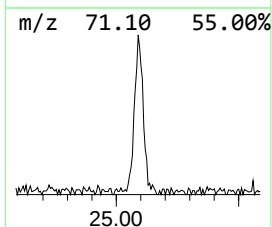
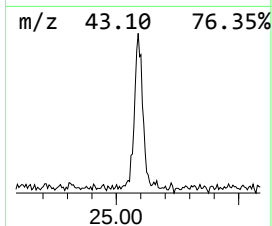
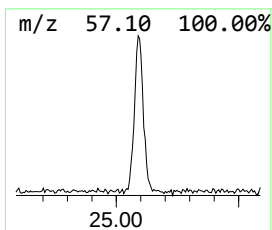
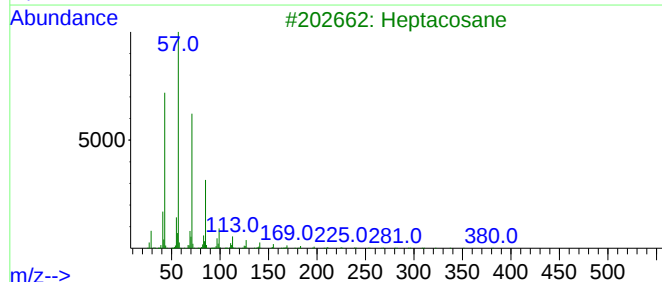
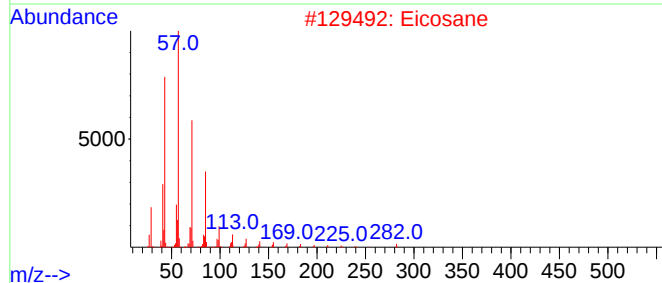
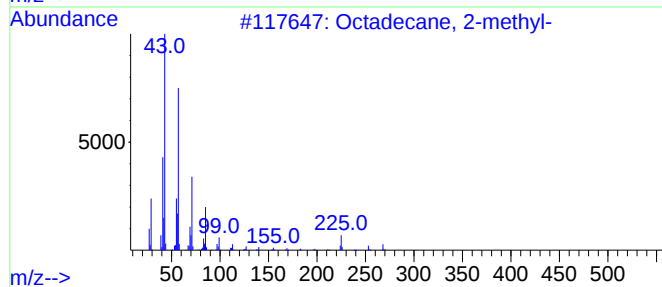
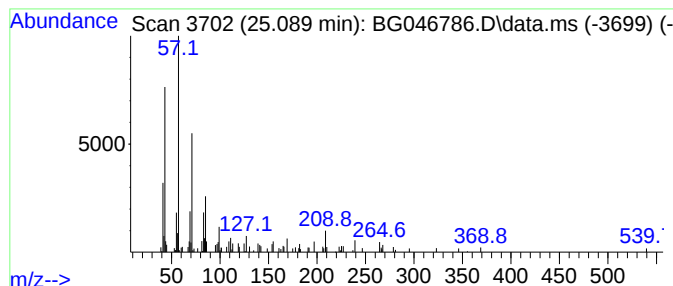
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.090	3.09 ng/ul	175302	Perylene-d12	25.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 2-methyl-	268	C19H40	001560-88-9	72
2	Eicosane	282	C20H42	000112-95-8	59
3	Heptacosane	380	C27H56	000593-49-7	59
4	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	59
5	Octadecane	254	C18H38	000593-45-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
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Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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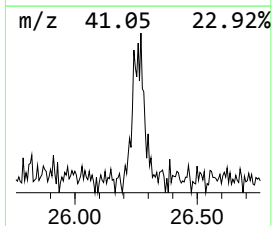
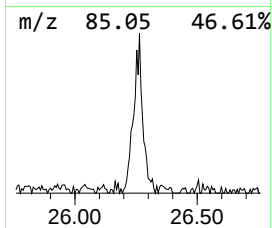
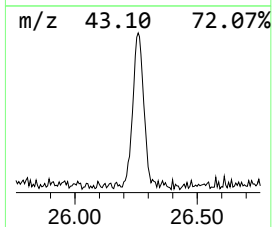
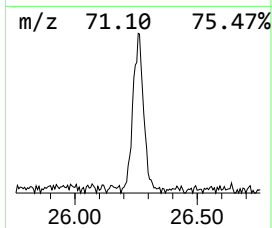
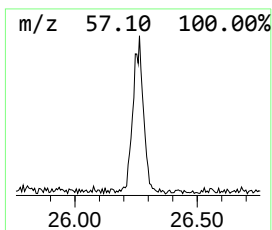
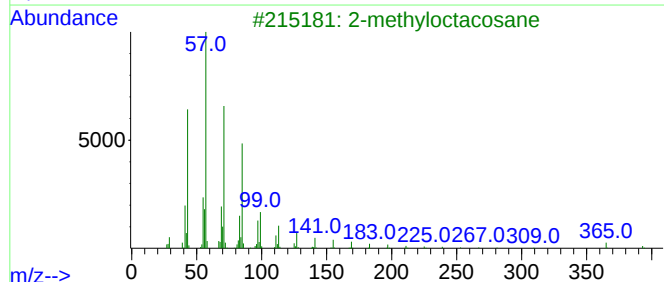
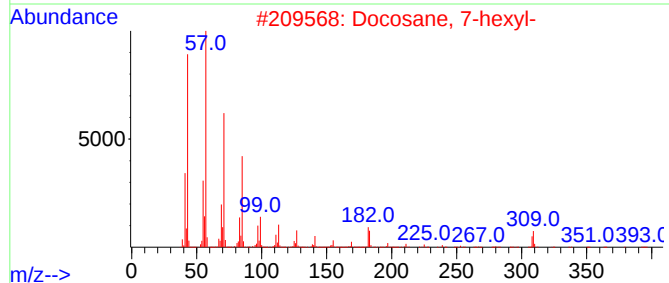
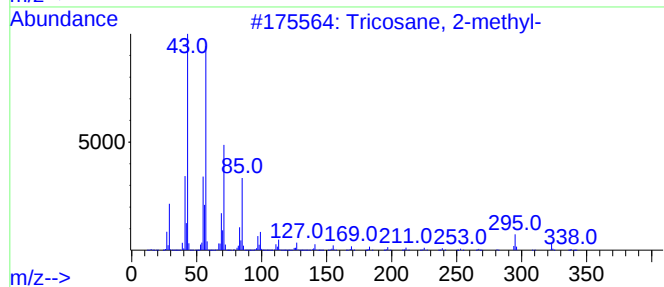
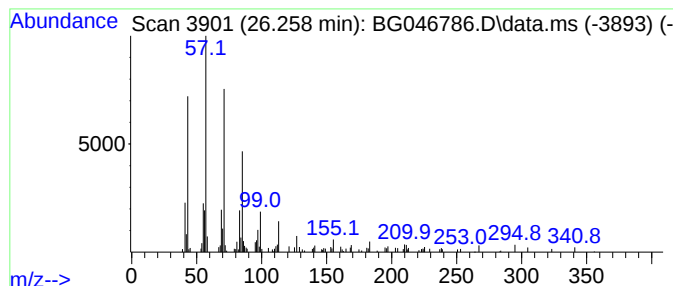
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.260	3.62 ng/ul	205298	Perylene-d12	25.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tricosane, 2-methyl-	338	C24H50	001928-30-9	91
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	86
4		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
Operator : CG/JU
Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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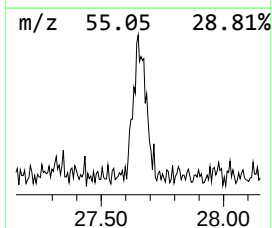
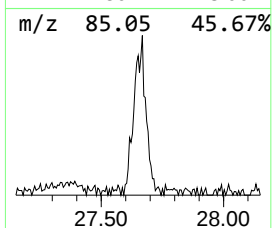
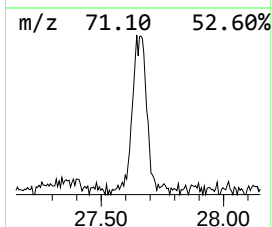
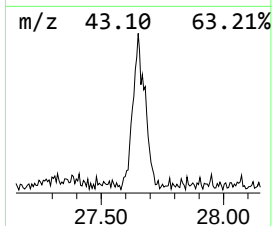
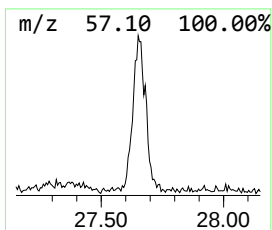
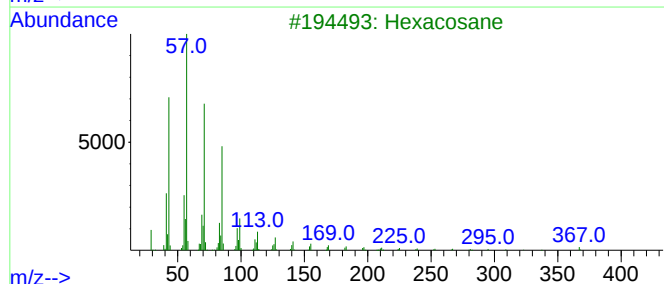
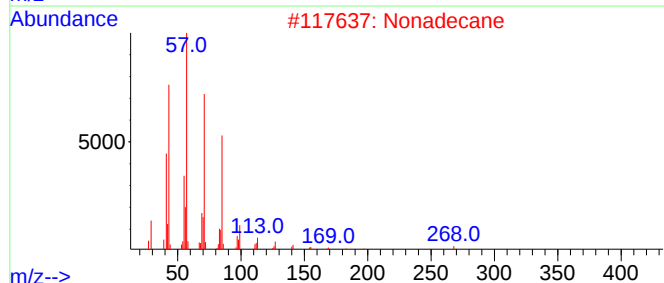
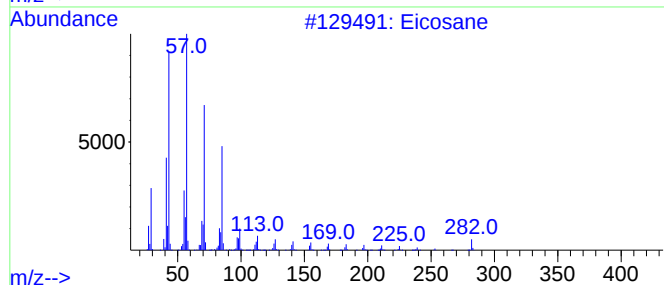
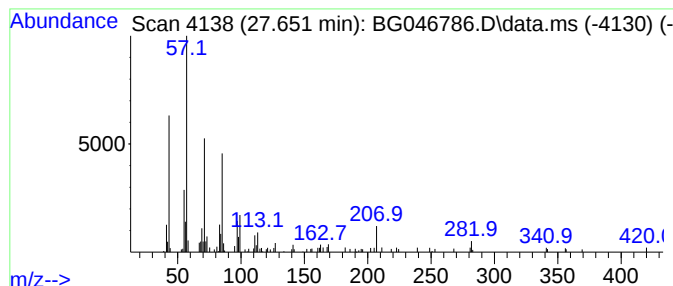
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.650	4.46 ng/ul	253028	Perylene-d12	25.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Nonadecane		268	C19H40	000629-92-5	86
3	Hexacosane		366	C26H54	000630-01-3	83
4	Octacosane		394	C28H58	000630-02-4	83
5	Octacosane		394	C28H58	000630-02-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
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Misc :
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Instrument :
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ClientSampleId :
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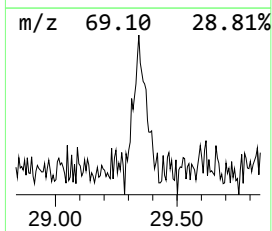
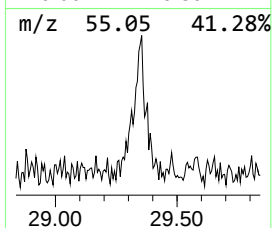
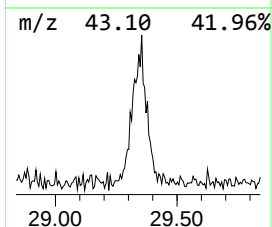
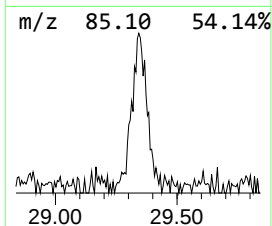
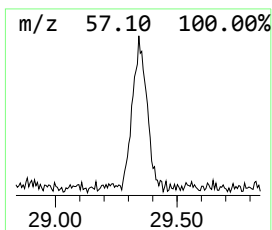
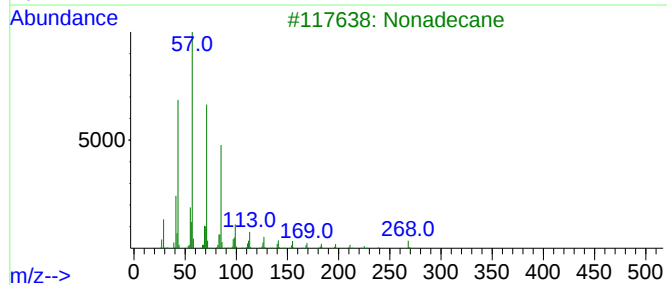
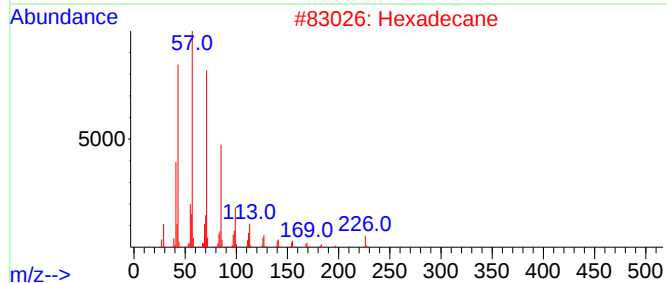
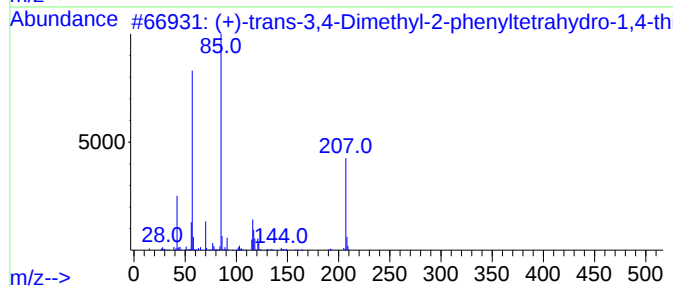
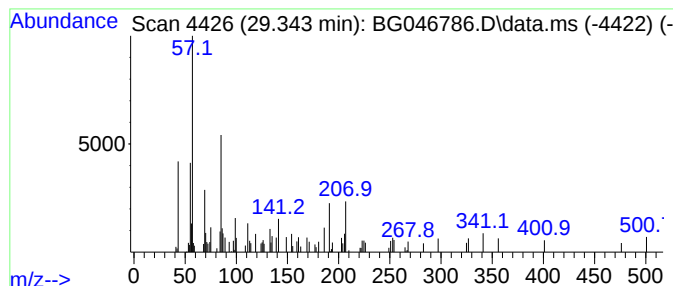
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.340	3.07 ng/ul	174235	Perylene-d12	25.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-trans-3,4-Dimethyl-2-phenylt...	207	C12H17NS	1000244-64-9	38		
2	Hexadecane	226	C16H34	000544-76-3	18		
3	Nonadecane	268	C19H40	000629-92-5	18		
4	Octadecane	254	C18H38	000593-45-3	18		
5	Propiohydrazide, 2,2-dimethyl-N2...	226	C12H22N2O2	1000264-71-8	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100520\
Data File : BG046786.D
Acq On : 5 Oct 2020 16:53
Operator : CG/JU
Sample : L4153-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM01-EPA-BG092820.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3,3-Dimethyl-4-...	16.700	3.4	ng/ul	163343	4	17.474	968662	20.0
1,1'-Biphenyl, ...	17.120	27.5	ng/ul	1332510	4	17.474	968662	20.0
Naphth[2,3-b]az...	20.890	3.4	ng/ul	204812	5	21.752	1210050	20.0
Benz(a)anthrace...	23.940	18.9	ng/ul	1073280	6	25.024	1135200	20.0
(DEL) Alkane: S...	24.110	2.4	ng/ul	137335	6	25.024	1135200	20.0
(DEL) Alkane: S...	25.090	3.1	ng/ul	175302	6	25.024	1135200	20.0
(DEL) Alkane: S...	26.260	3.6	ng/ul	205298	6	25.024	1135200	20.0
(DEL) Alkane: S...	27.650	4.5	ng/ul	253028	6	25.024	1135200	20.0
unknown-01	29.340	3.1	ng/ul	174235	6	25.024	1135200	20.0