

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
 Data File : BG052769.D
 Acq On : 21 Mar 2022 15:52
 Operator : CG/JU
 Sample : N1708-18 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PPSP-B10-25.5-26.0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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\8270-BG031522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG052769.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.783	569	578	591	rVB4	70787	210102	10.75%	0.412%
2	7.136	628	638	644	rBV3	111090	248741	12.73%	0.488%
3	7.294	659	665	673	rVB4	183192	369378	18.91%	0.725%
4	7.794	744	750	758	rVV	317798	548351	28.07%	1.076%
5	8.058	786	795	802	rBV4	64327	168630	8.63%	0.331%
6	8.140	802	809	816	rVB2	277995	611626	31.31%	1.200%
7	8.269	826	831	839	rVB2	119414	239643	12.27%	0.470%
8	8.446	857	861	867	rVB2	95841	166587	8.53%	0.327%
9	8.798	915	921	927	rBV3	141506	354448	18.14%	0.695%
10	8.921	935	942	945	rBV2	102474	192812	9.87%	0.378%
11	9.156	977	982	990	rVB	137919	265596	13.60%	0.521%
12	9.262	990	1000	1005	rBV3	284177	649333	33.24%	1.274%
13	9.344	1009	1014	1024	rVB4	113527	282438	14.46%	0.554%
14	9.503	1032	1041	1048	rBV10	92260	343646	17.59%	0.674%
15	9.591	1051	1056	1061	rBV4	76638	189266	9.69%	0.371%
16	9.791	1084	1090	1095	rVV3	115829	265759	13.60%	0.521%
17	9.844	1095	1099	1103	rVV	151480	253134	12.96%	0.497%
18	10.043	1124	1133	1137	rBV3	116687	266924	13.66%	0.524%
19	10.090	1137	1141	1145	rVB3	109862	181256	9.28%	0.356%
20	10.143	1145	1150	1156	rBV6	155312	281644	14.42%	0.553%
21	10.208	1156	1161	1164	rBV3	153629	274021	14.03%	0.538%
22	10.302	1171	1177	1181	rBV2	112143	197085	10.09%	0.387%
23	10.355	1181	1186	1193	rVV3	268326	620472	31.76%	1.217%
24	10.419	1193	1197	1203	rVB4	161885	303244	15.52%	0.595%
25	10.490	1203	1209	1212	rBV	131888	247395	12.66%	0.485%
26	10.537	1214	1217	1222	rVV3	107208	189535	9.70%	0.372%
27	10.590	1222	1226	1233	rVB	189841	341760	17.49%	0.671%
28	10.660	1233	1238	1242	rBV	99705	180032	9.22%	0.353%
29	10.954	1278	1288	1293	rVV2	386638	1173437	60.07%	2.302%
30	11.007	1293	1297	1305	rVV2	353315	876960	44.89%	1.721%
31	11.077	1305	1309	1312	rVV	171381	321326	16.45%	0.630%
32	11.124	1312	1317	1322	rVV	351746	606071	31.02%	1.189%
33	11.183	1322	1327	1334	rVV3	112903	253153	12.96%	0.497%
34	11.430	1364	1369	1380	rVB3	245032	575446	29.46%	1.129%
35	11.541	1381	1388	1397	rBV5	133764	383273	19.62%	0.752%
36	11.624	1398	1402	1405	rVV2	146688	257348	13.17%	0.505%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
 Data File : BG052769.D
 Acq On : 21 Mar 2022 15:52
 Operator : CG/JU
 Sample : N1708-18 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PPSP-B10-25.5-26.0

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % 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\8270-BG031522.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.665	1405	1409	1416	rVV5	218446	515296	26.38%	1.011%
38	11.723	1416	1419	1426	rVB3	112611	225547	11.55%	0.443%
39	11.794	1426	1431	1437	rBV3	175624	299269	15.32%	0.587%
40	11.876	1438	1445	1451	rBV4	229268	504123	25.81%	0.989%
41	11.982	1457	1463	1467	rBV	426359	710084	36.35%	1.393%
42	12.064	1473	1477	1483	rVB2	162246	279632	14.31%	0.549%
43	12.146	1485	1491	1498	rVB2	550692	996714	51.02%	1.955%
44	12.499	1546	1551	1556	rVB3	306150	523622	26.80%	1.027%
45	12.605	1560	1569	1575	rVB2	934030	1951026	99.87%	3.828%
46	12.699	1580	1585	1592	rBV6	61988	165363	8.46%	0.324%
47	12.822	1599	1606	1609	rBV	573638	1111352	56.89%	2.180%
48	12.857	1609	1612	1617	rVB2	417436	630278	32.26%	1.237%
49	12.998	1631	1636	1639	rBV3	178074	277707	14.22%	0.545%
50	13.039	1639	1643	1645	rBV4	120950	194383	9.95%	0.381%
51	13.104	1650	1654	1661	rVV4	178765	340387	17.42%	0.668%
52	13.263	1677	1681	1685	rBV5	96216	212073	10.86%	0.416%
53	13.315	1685	1690	1696	rVB2	607981	960273	49.16%	1.884%
54	13.603	1730	1739	1745	rBV6	200579	567689	29.06%	1.114%
55	13.715	1754	1758	1763	rVB2	304599	513996	26.31%	1.008%
56	13.768	1763	1767	1769	rBV3	161894	227383	11.64%	0.446%
57	13.803	1769	1773	1776	rBV	189287	248766	12.73%	0.488%
58	13.932	1789	1795	1810	rVB3	625014	1836988	94.03%	3.604%
59	14.091	1811	1822	1827	rBV2	878400	1953542	100.00%	3.833%
60	14.144	1827	1831	1837	rVB2	664323	1113521	57.00%	2.185%
61	14.202	1837	1841	1843	rBV2	158786	235616	12.06%	0.462%
62	14.249	1845	1849	1853	rVB	564098	759126	38.86%	1.489%
63	14.326	1858	1862	1872	rVB4	392733	1017815	52.10%	1.997%
64	14.490	1887	1890	1894	rVB	130764	172805	8.85%	0.339%
65	14.608	1906	1910	1915	rBV4	200466	294106	15.06%	0.577%
66	14.737	1927	1932	1934	rBV	415187	675712	34.59%	1.326%
67	14.766	1934	1937	1942	rVB2	457110	677364	34.67%	1.329%
68	14.819	1942	1946	1948	rBV	102485	168150	8.61%	0.330%
69	14.972	1967	1972	1981	rVB	387645	1010561	51.73%	1.983%
70	15.054	1982	1986	1990	rBV2	117268	202787	10.38%	0.398%
71	15.160	2000	2004	2012	rVB	561132	940414	48.14%	1.845%
72	15.236	2012	2017	2021	rVB	419638	606315	31.04%	1.190%
73	15.283	2021	2025	2036	rVB7	261363	624481	31.97%	1.225%
74	15.383	2036	2042	2046	rBV	364508	735563	37.65%	1.443%
75	15.436	2047	2051	2055	rVB	346359	504804	25.84%	0.990%
76	15.671	2085	2091	2094	rBV5	101286	240531	12.31%	0.472%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
 Data File : BG052769.D
 Acq On : 21 Mar 2022 15:52
 Operator : CG/JU
 Sample : N1708-18 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PPSP-B10-25.5-26.0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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\8270-BG031522.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.718	2095	2099	2106	rVB4	195903	404671	20.71%	0.794%
78	15.800	2109	2113	2117	rBV2	221195	340756	17.44%	0.669%
79	15.853	2117	2122	2126	rVB3	157043	330445	16.92%	0.648%
80	15.935	2133	2136	2139	rVB3	195280	241596	12.37%	0.474%
81	15.977	2139	2143	2146	rVV4	145119	242801	12.43%	0.476%
82	16.018	2146	2150	2157	rVB2	715976	1071489	54.85%	2.102%
83	16.223	2182	2185	2189	rVV2	131871	193085	9.88%	0.379%
84	16.317	2198	2201	2206	rBV3	110936	191540	9.80%	0.376%
85	16.494	2226	2231	2242	rVB2	1114980	1868069	95.62%	3.665%
86	16.869	2289	2295	2300	rBV4	386735	719016	36.81%	1.411%
87	17.316	2366	2371	2382	rVB2	746416	1420651	72.72%	2.787%
88	17.510	2399	2404	2408	rBV	522654	834990	42.74%	1.638%
89	17.551	2408	2411	2415	rVB	251357	350999	17.97%	0.689%
90	17.704	2434	2437	2442	rVB5	111826	169481	8.68%	0.333%
91	17.921	2470	2474	2485	rVB4	222652	429561	21.99%	0.843%
92	18.385	2549	2553	2557	rBV	299780	442698	22.66%	0.869%
93	18.432	2557	2561	2568	rVB	393145	624396	31.96%	1.225%
94	18.567	2581	2584	2588	rVB2	137003	189857	9.72%	0.372%
95	19.296	2703	2708	2712	rVV	268808	397542	20.35%	0.780%
96	19.930	2812	2816	2822	rVB3	110719	196666	10.07%	0.386%
97	19.989	2822	2826	2832	rVB	129602	193474	9.90%	0.380%
98	20.106	2843	2846	2851	rVB	156597	188723	9.66%	0.370%
99	21.792	3126	3133	3142	rVB	507194	879444	45.02%	1.725%
100	25.111	3688	3698	3712	rVB2	315991	933423	47.78%	1.831%

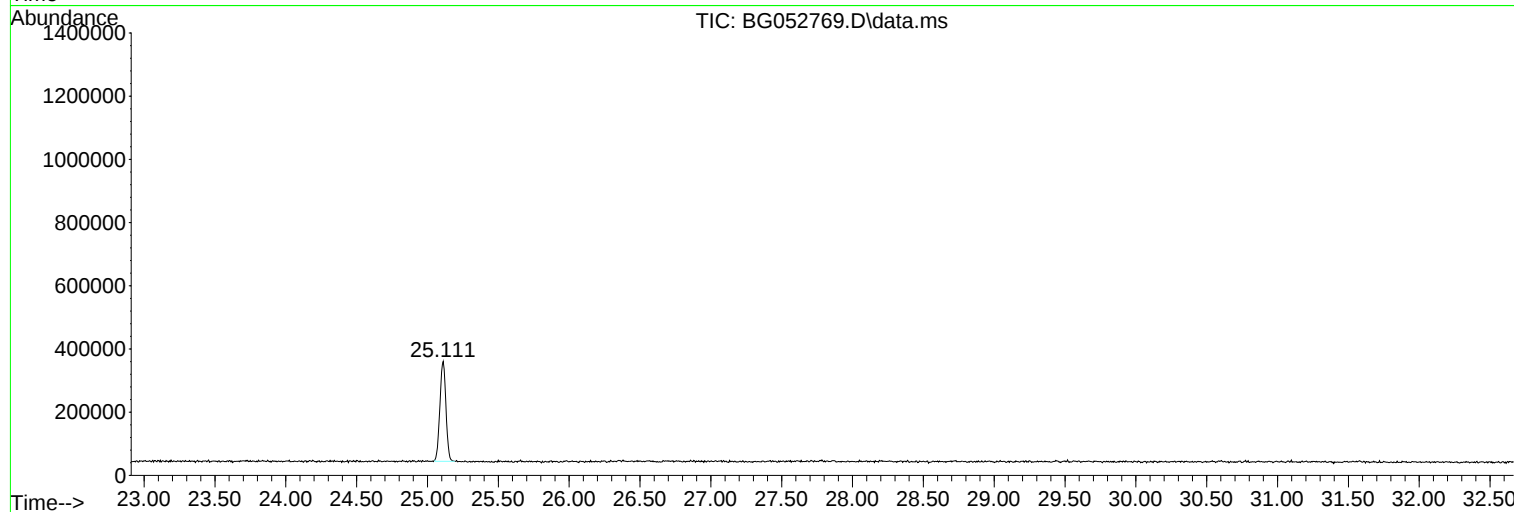
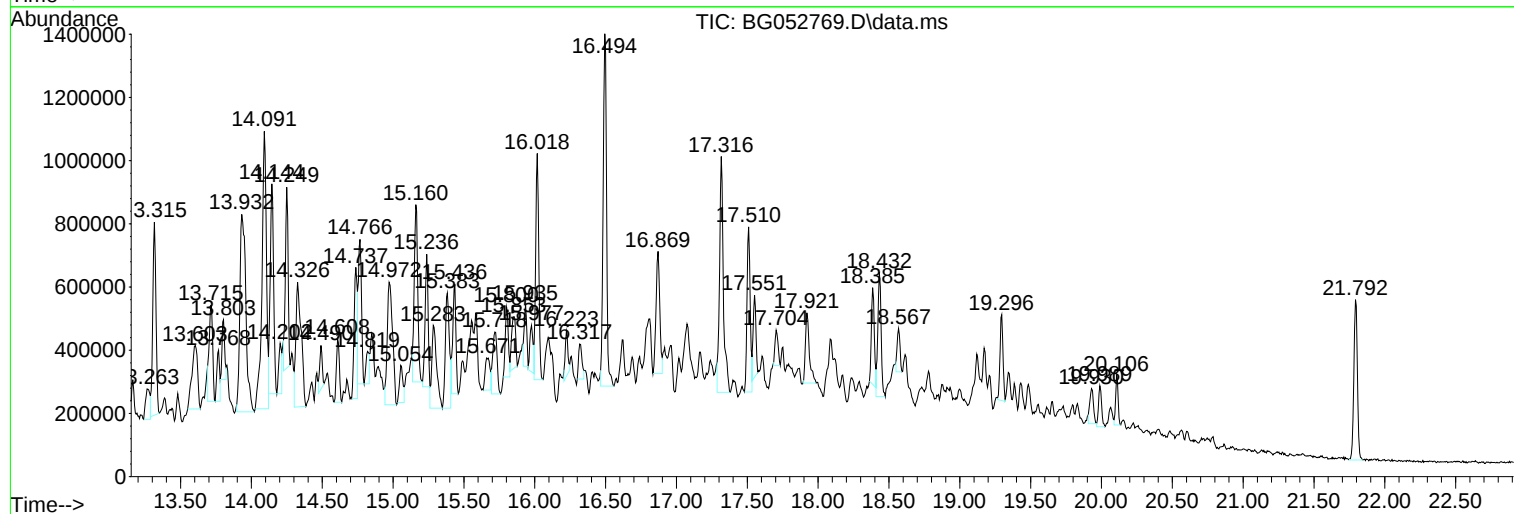
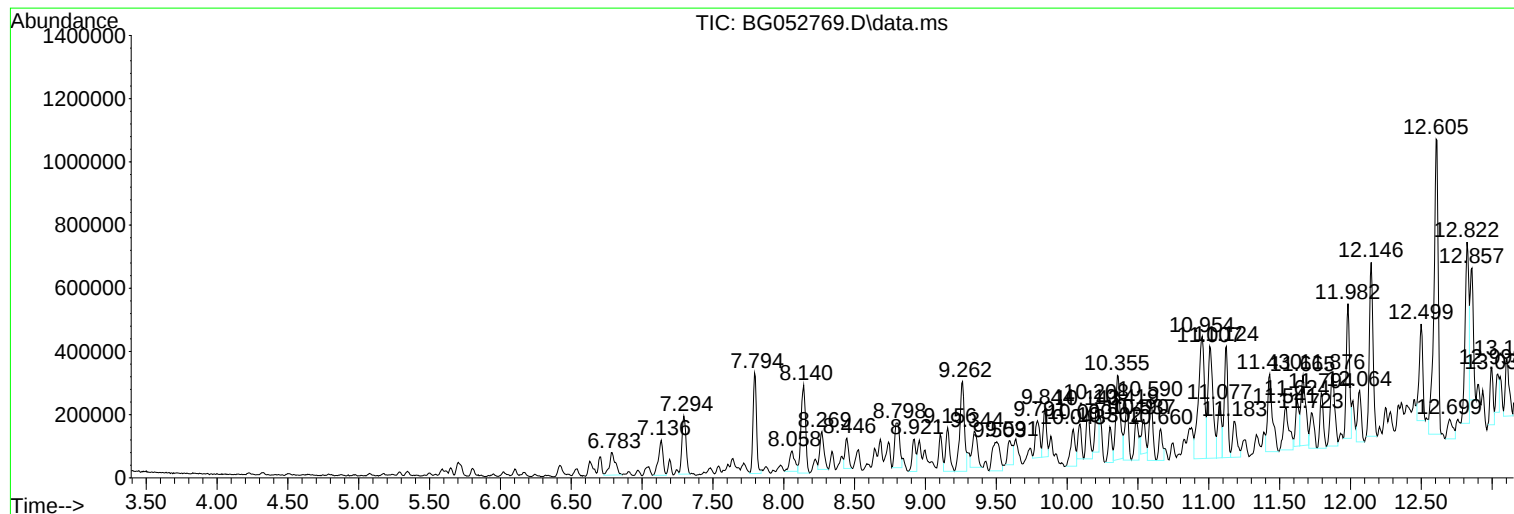
Sum of corrected areas: 50970409

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

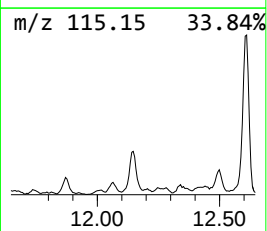
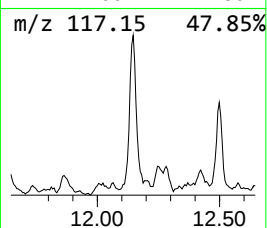
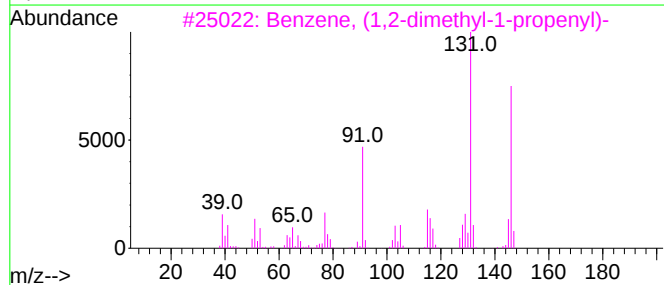
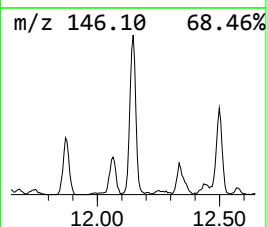
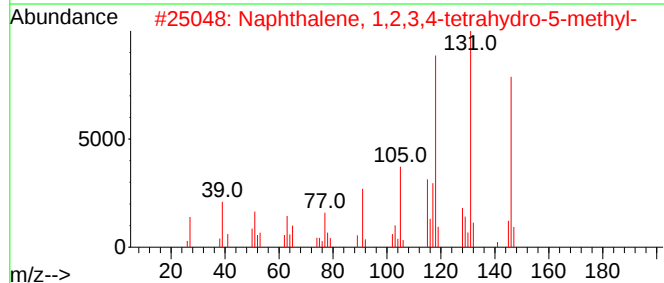
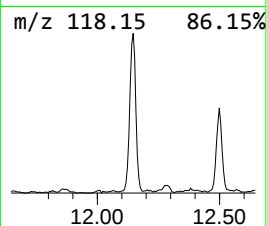
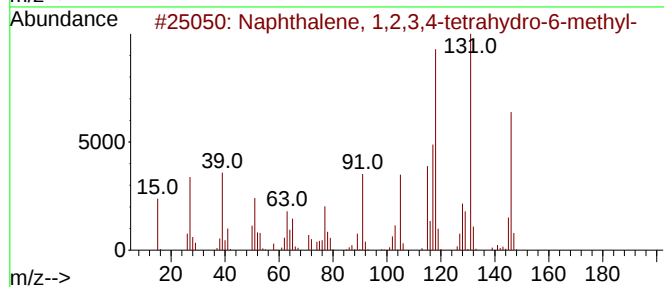
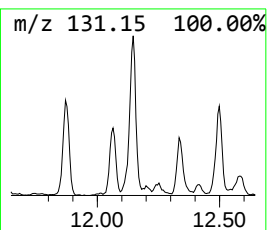
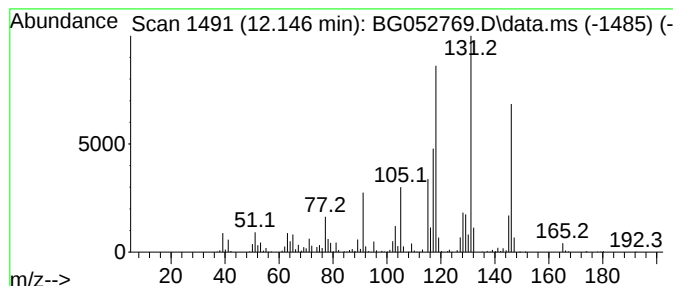
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.146	16.99 ng	996714	Naphthalene-d8	10.960	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
4	Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	74
5	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

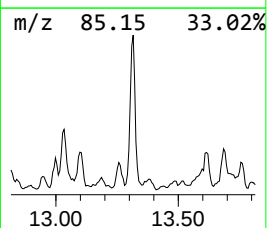
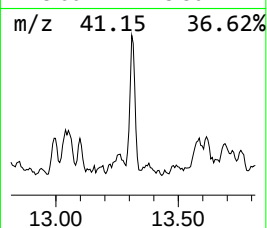
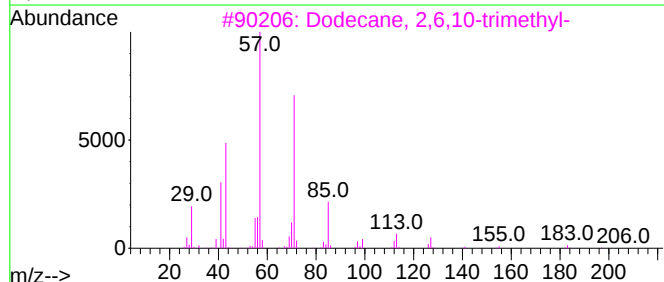
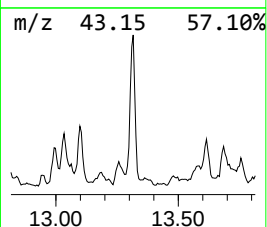
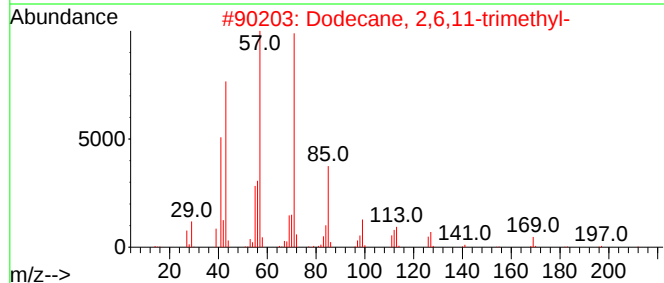
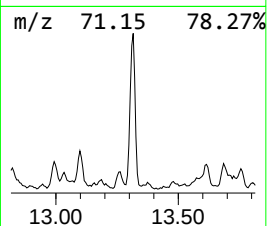
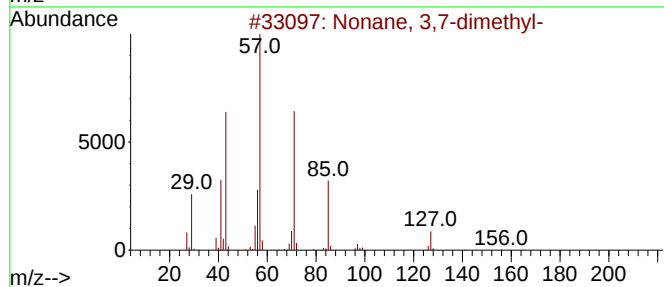
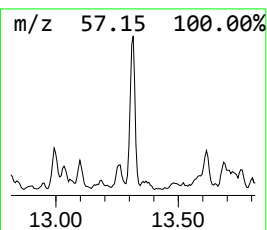
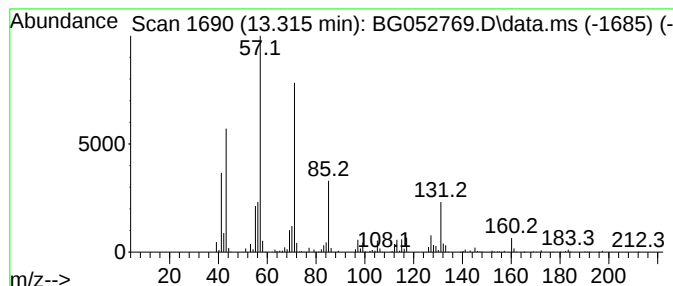
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Nonane, 3,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.316	28.35 ng	960273	Acenaphthene-d10	14.766

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	64
2			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	58
4			Decane, 5-propyl-	184	C13H28	017312-62-8	53
5			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

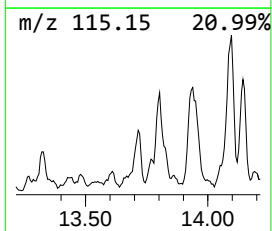
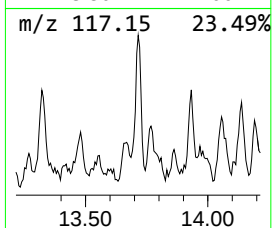
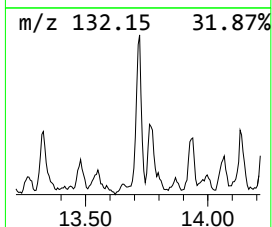
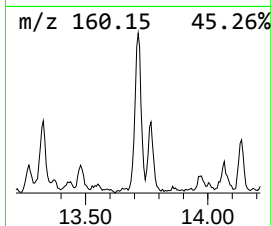
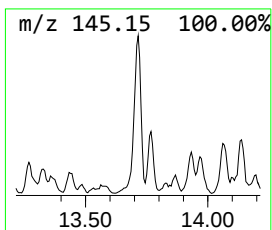
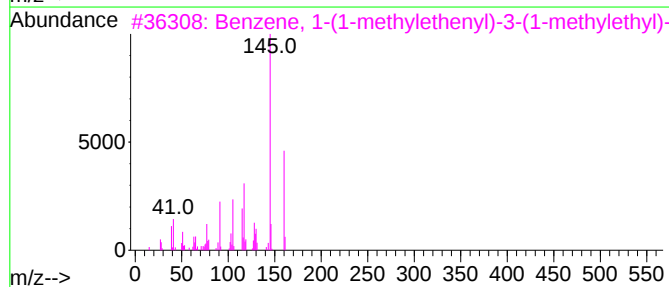
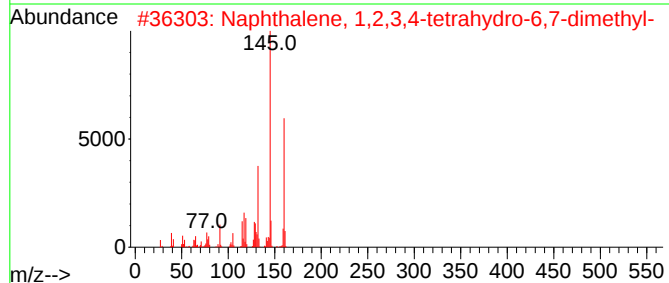
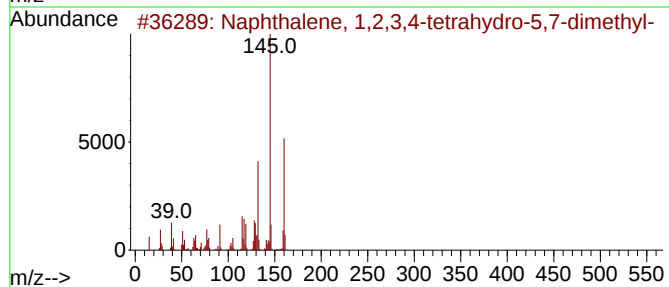
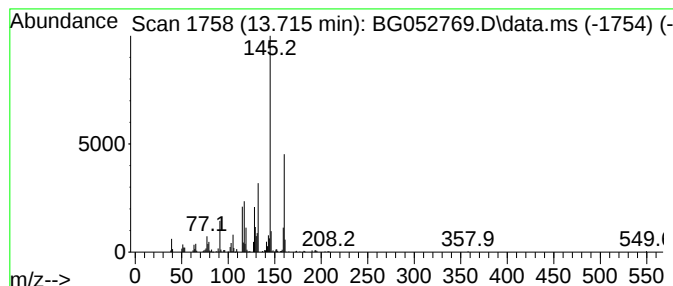
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.715	15.18 ng	513996	Acenaphthene-d10	14.766	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	94
3	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	89
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	020027-77-4	87
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

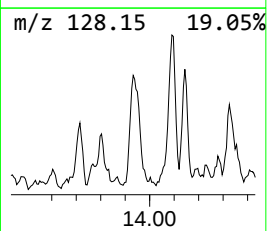
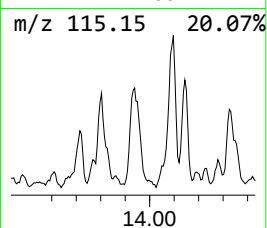
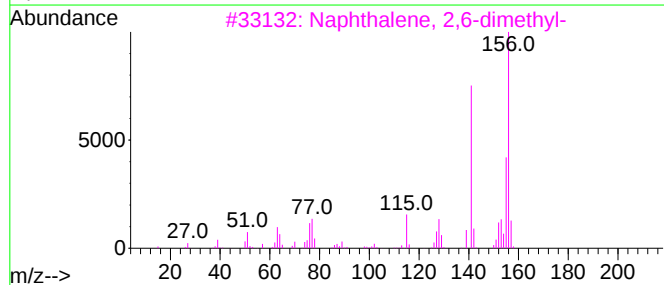
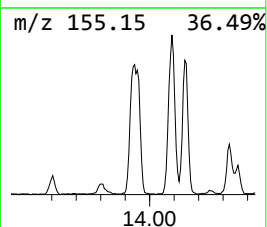
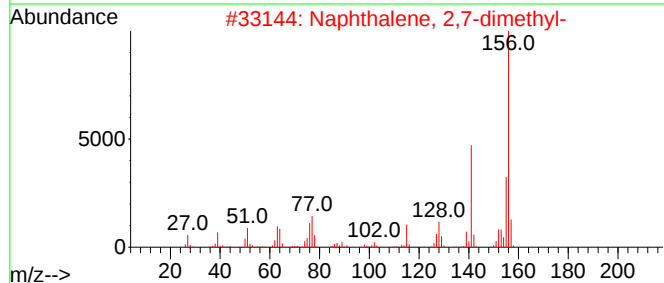
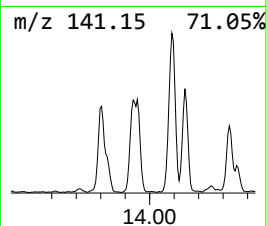
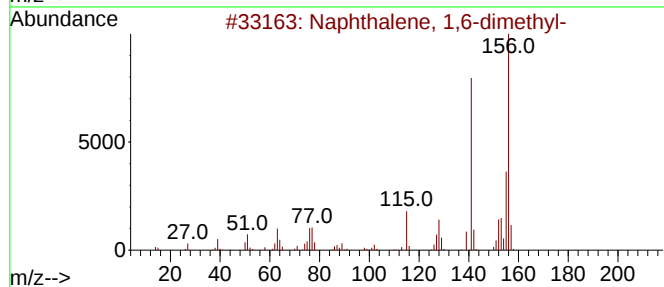
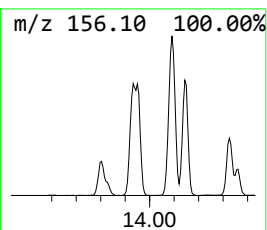
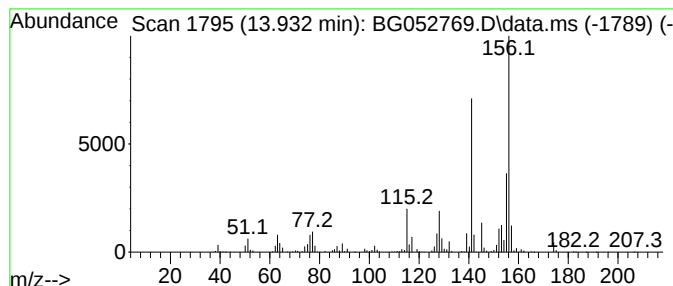
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.932	54.24 ng	1836990	Acenaphthene-d10	14.766

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

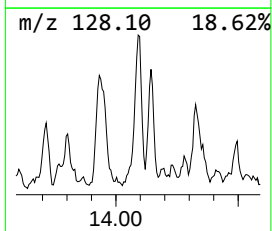
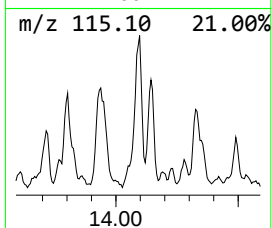
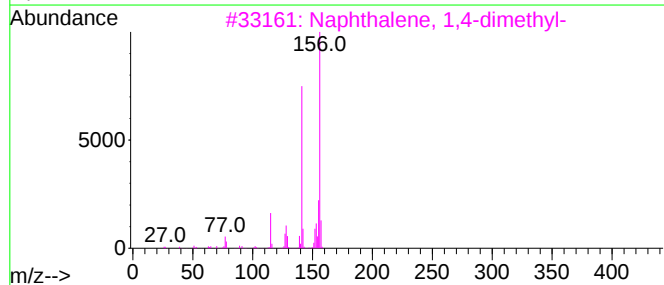
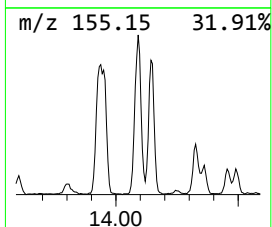
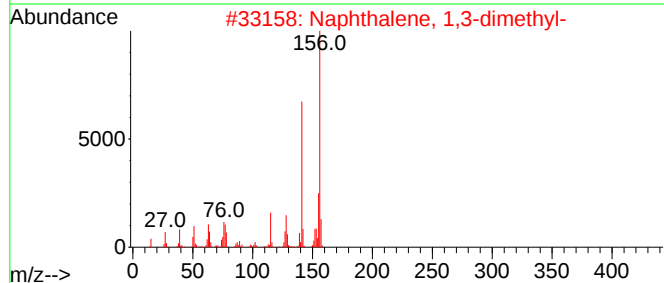
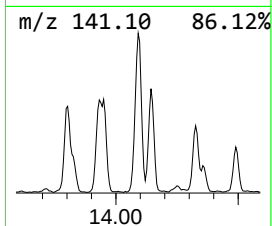
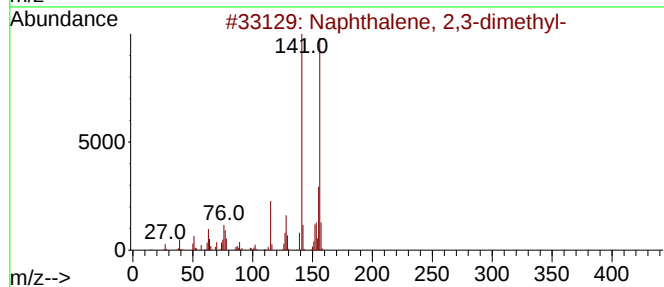
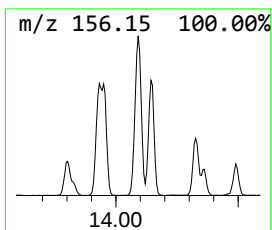
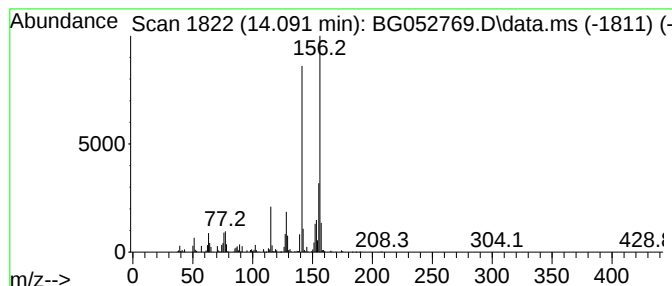
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.091	57.68 ng	1953540	Acenaphthene-d10		14.766	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97
2	Naphthalene, 1,3-dimethyl-		156	C12H12	000575-41-7	97
3	Naphthalene, 1,4-dimethyl-		156	C12H12	000571-58-4	97
4	Naphthalene, 1,7-dimethyl-		156	C12H12	000575-37-1	96
5	Naphthalene, 1,6-dimethyl-		156	C12H12	000575-43-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

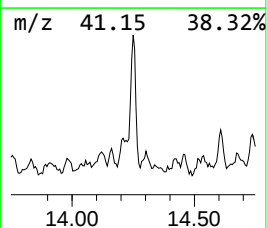
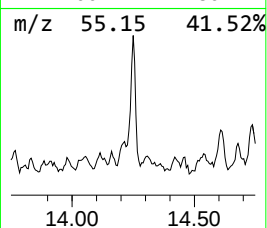
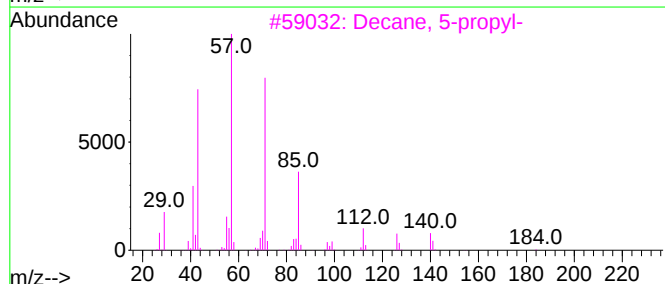
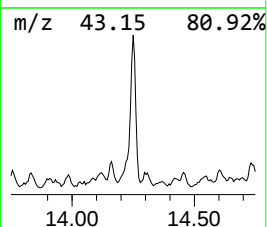
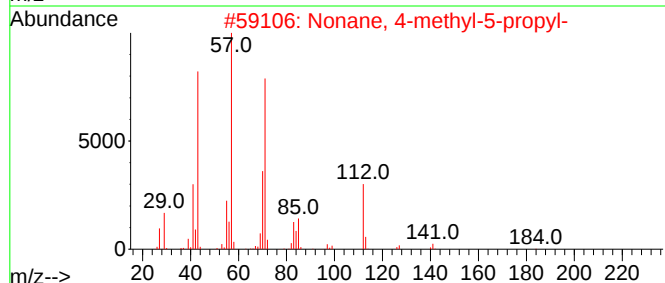
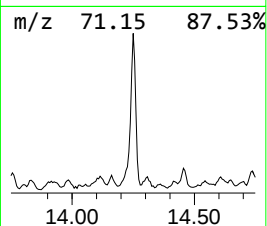
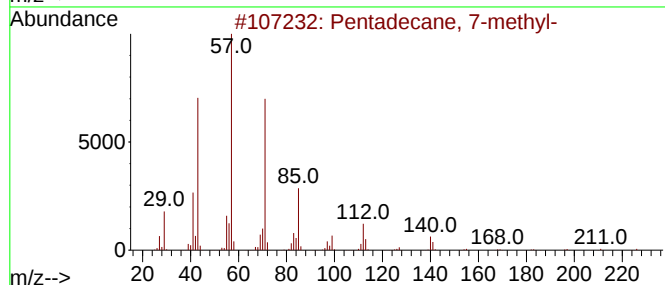
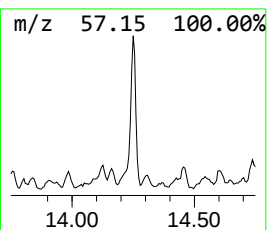
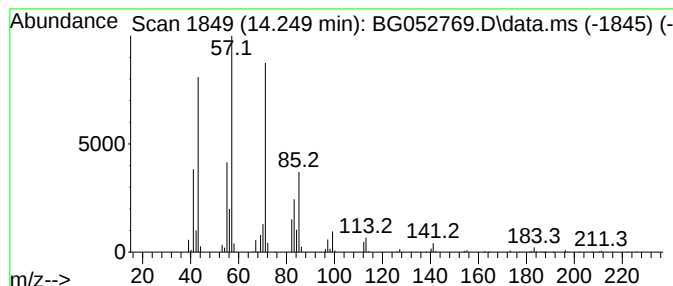
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 7-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.249	22.41 ng	759126	Acenaphthene-d10	14.766

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72
2			Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	68
3			Decane, 5-propyl-	184	C13H28	017312-62-8	68
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64
5			Tridecane, 5-propyl-	226	C16H34	055045-11-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

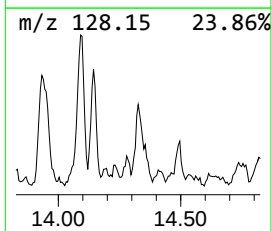
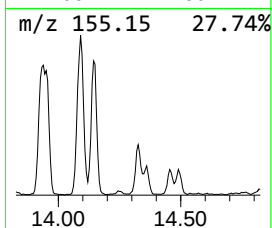
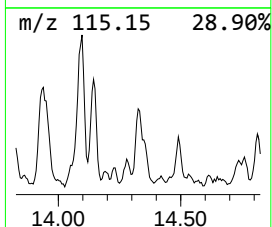
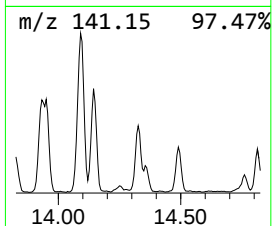
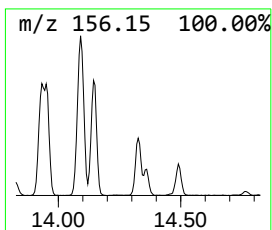
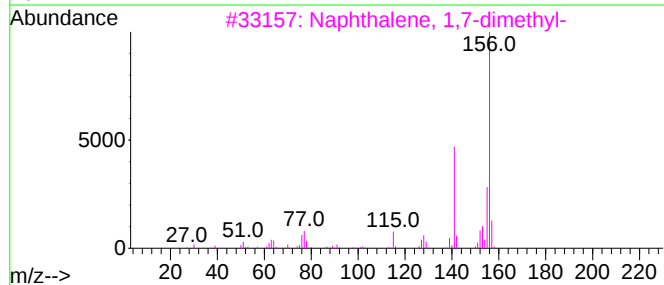
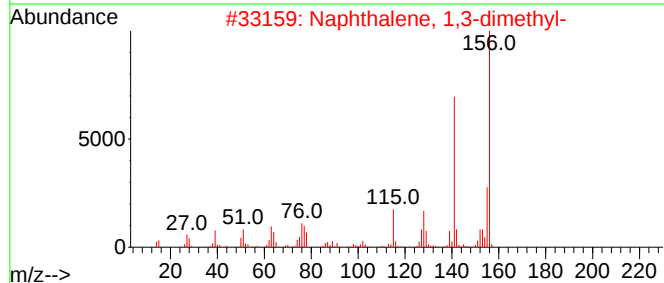
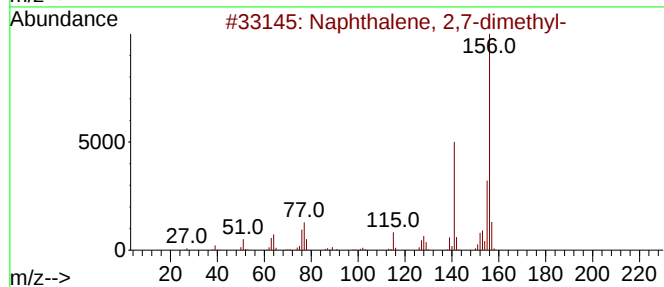
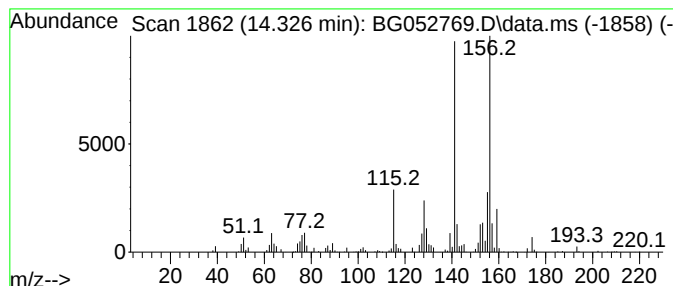
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.326	30.05 ng	1017820	Acenaphthene-d10	14.766

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

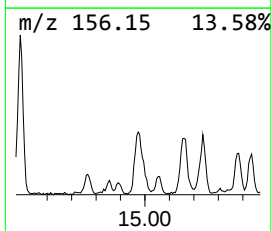
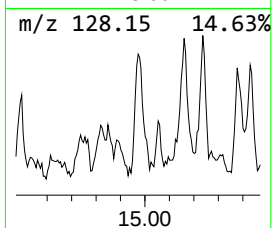
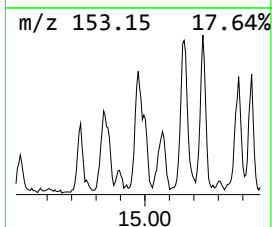
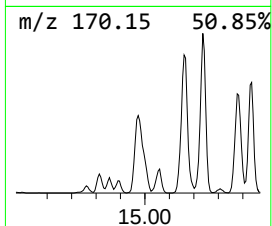
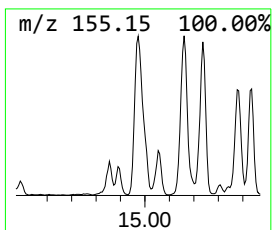
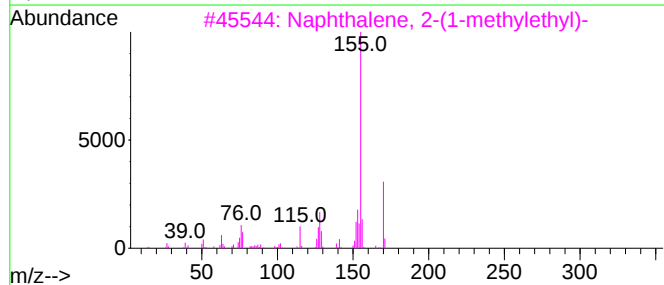
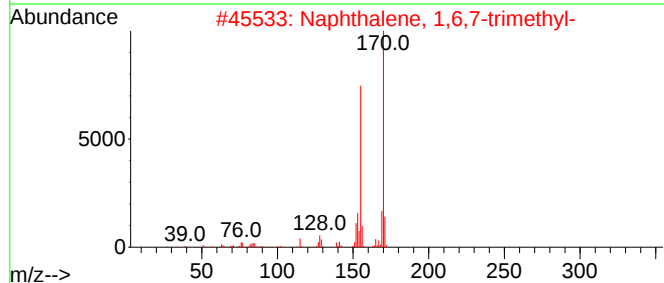
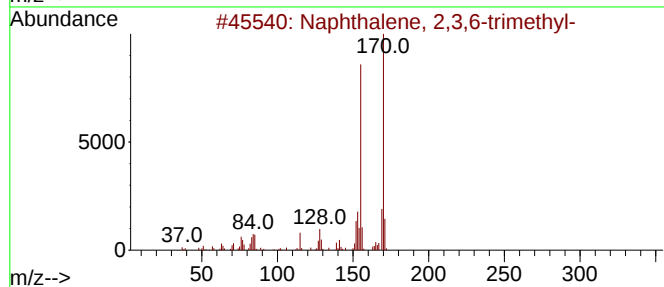
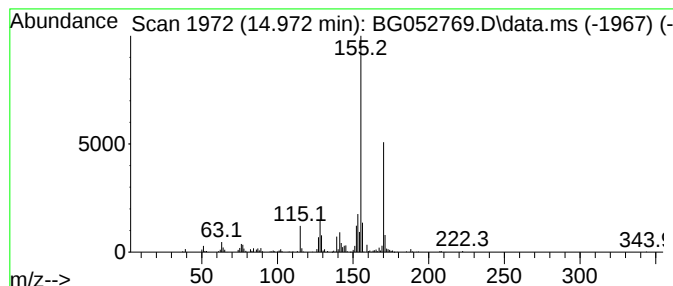
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,3,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.972	29.84 ng	1010560	Acenaphthene-d10	14.766	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
3	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	93
4	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	91
5	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

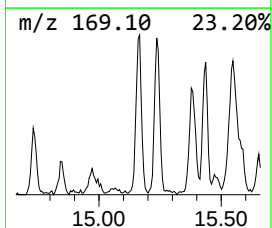
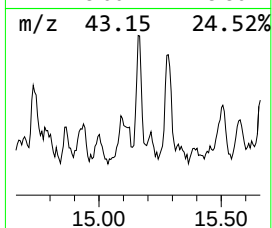
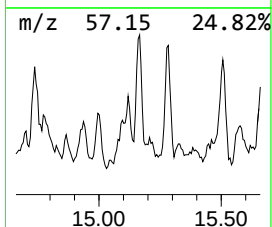
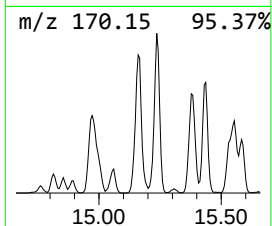
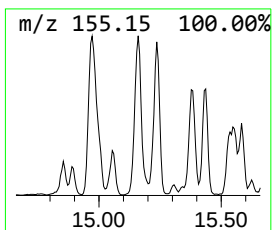
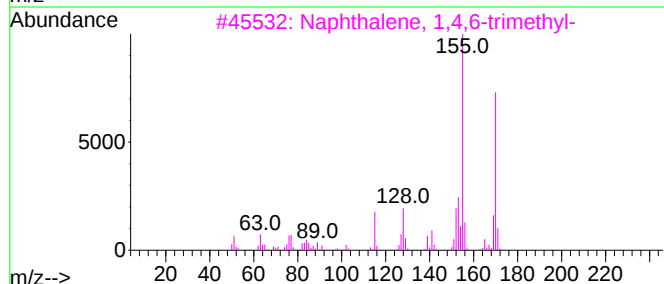
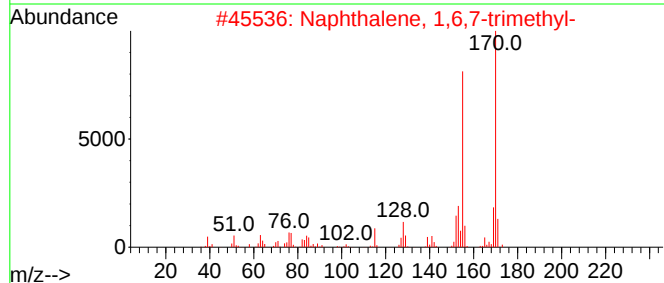
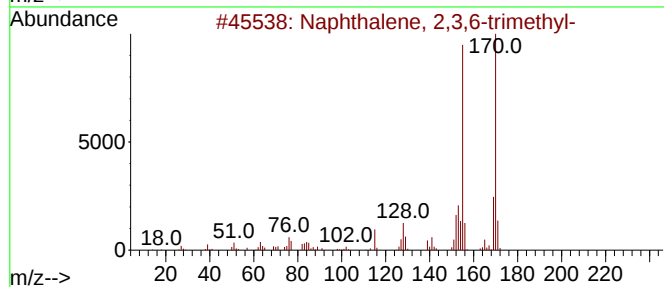
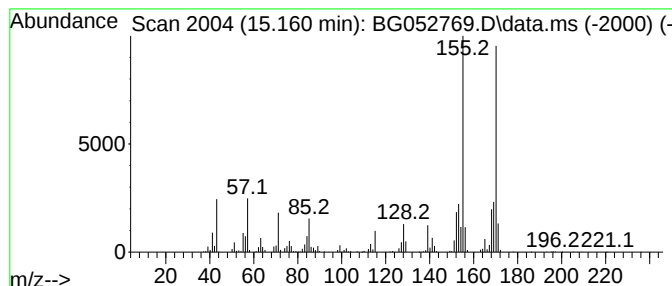
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,6,7-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.160	27.77 ng	940414	Acenaphthene-d10	14.766

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
4		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	90
5		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

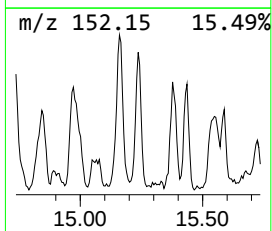
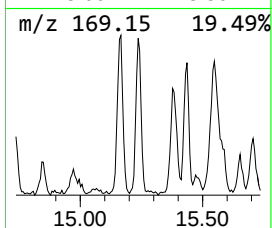
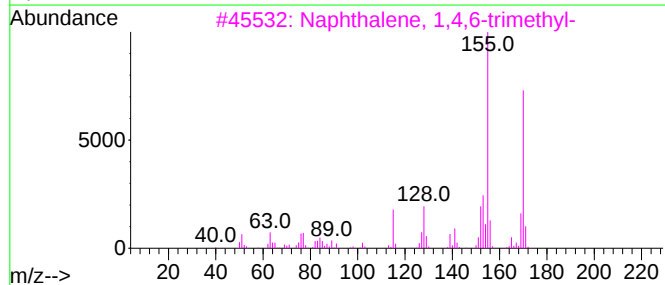
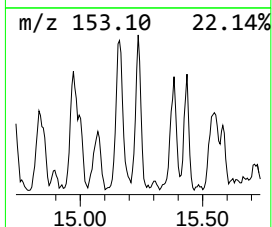
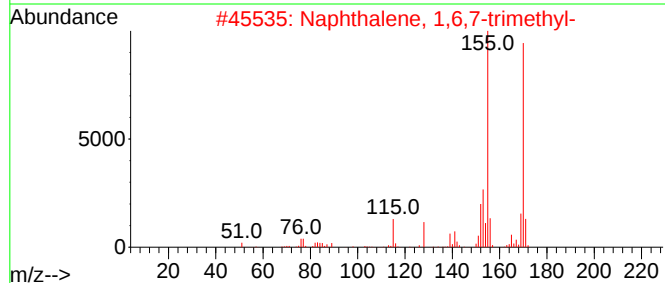
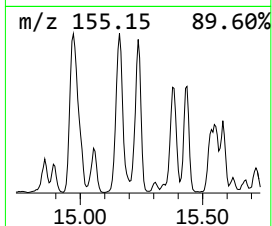
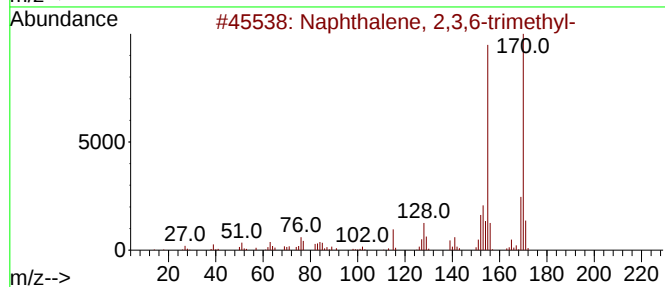
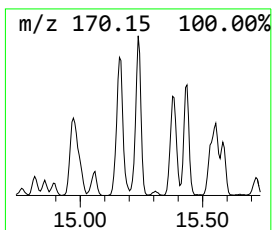
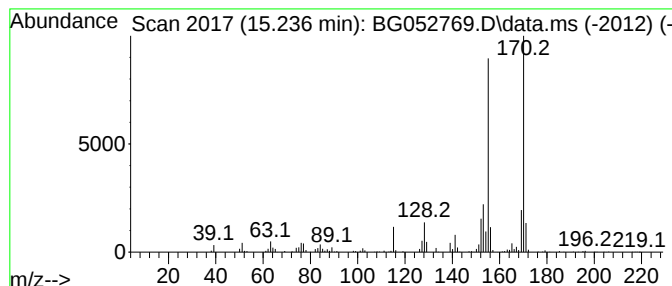
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1,4,6-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.236	17.90 ng	606315	Acenaphthene-d10		14.766	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-		170	C13H14	000829-26-5	97
2	Naphthalene, 1,6,7-trimethyl-		170	C13H14	002245-38-7	96
3	Naphthalene, 1,4,6-trimethyl-		170	C13H14	002131-42-2	96
4	1-(2,2-Dimethylcyclopropyl)-2-ph...		170	C13H14	1000294-91-1	94
5	Naphthalene, 1,4,5-trimethyl-		170	C13H14	002131-41-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

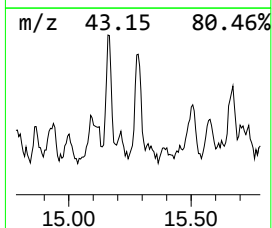
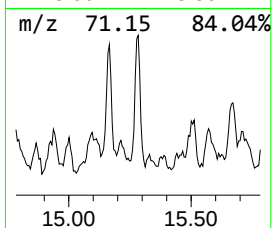
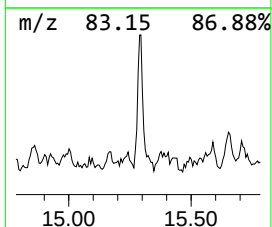
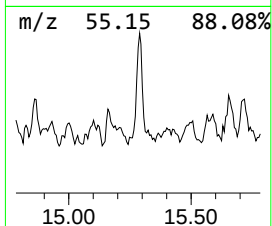
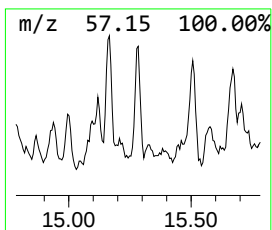
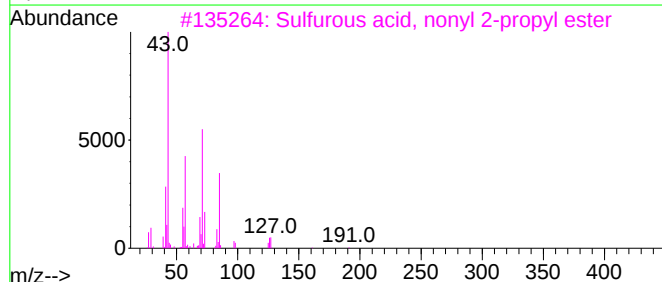
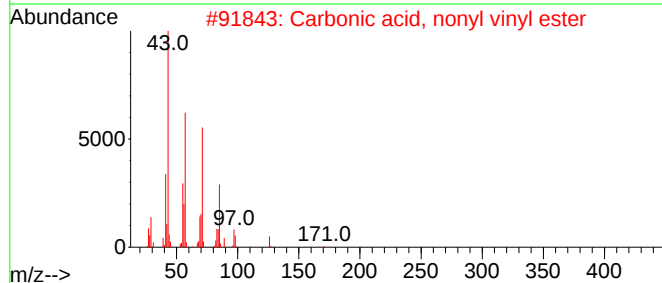
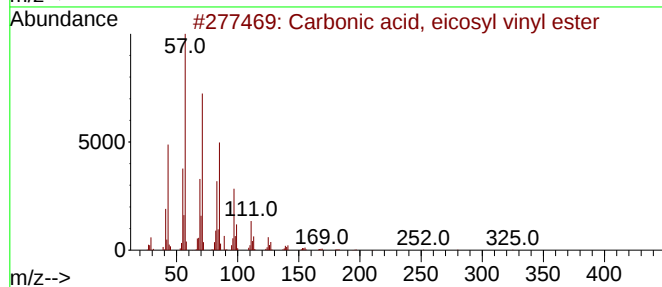
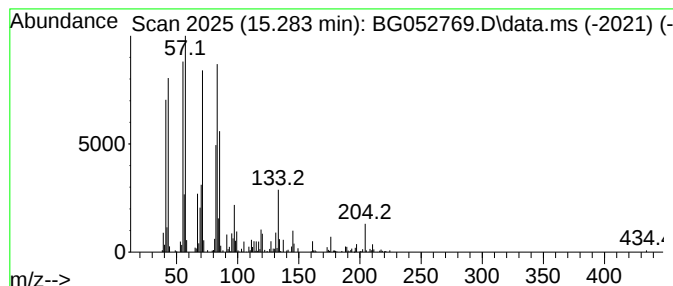
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.283 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.283	18.44 ng	624481	Acenaphthene-d10	14.766

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	38
2			Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	38
3			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	27
4			Sulfurous acid, decyl 2-propyl e...	264	C13H28O3S	1000309-12-1	27
5			1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

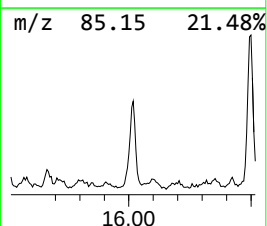
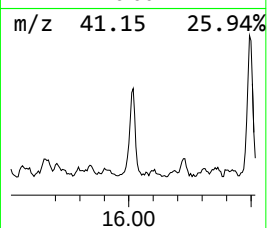
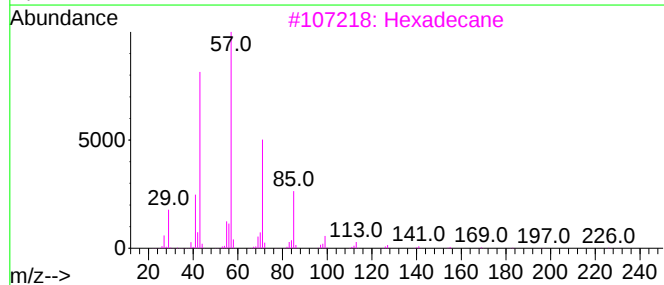
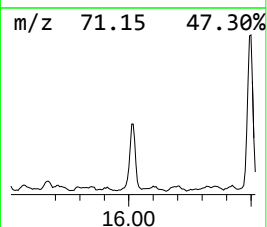
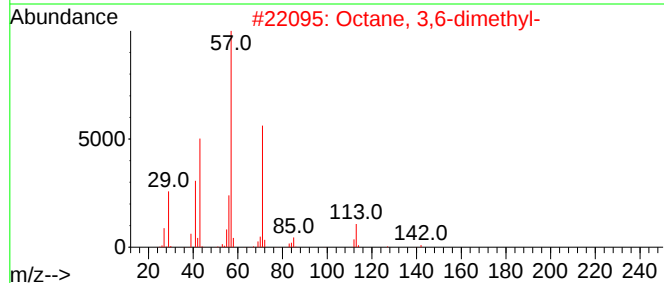
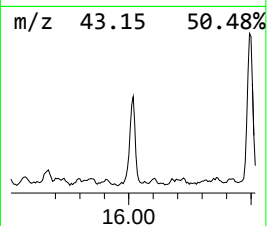
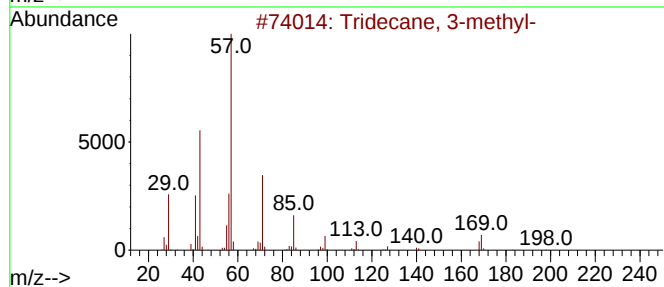
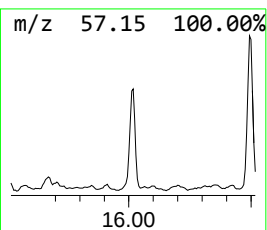
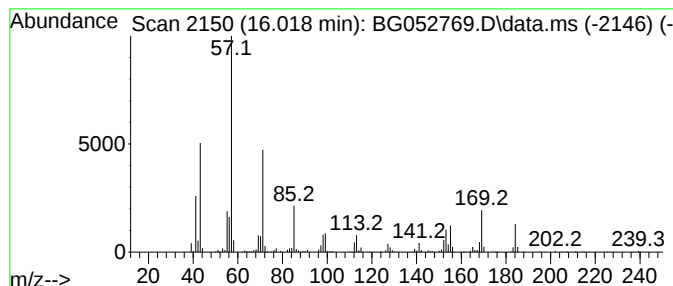
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tridecane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.018	31.64 ng	1071490	Acenaphthene-d10		14.766	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 3-methyl-		198	C14H30	006418-41-3	50
2	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	46
3	Hexadecane		226	C16H34	000544-76-3	46
4	Octane, 2-methyl-		128	C9H20	003221-61-2	46
5	Hexadecane, 3-methyl-		240	C17H36	006418-43-5	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

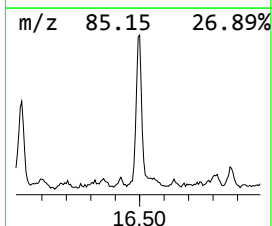
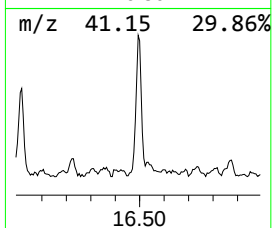
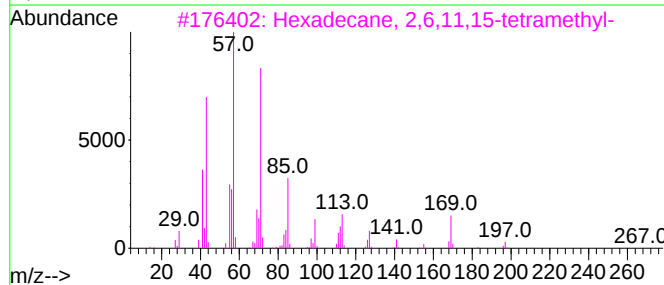
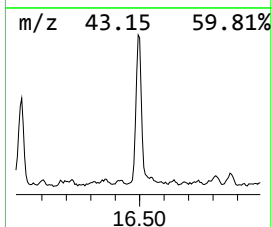
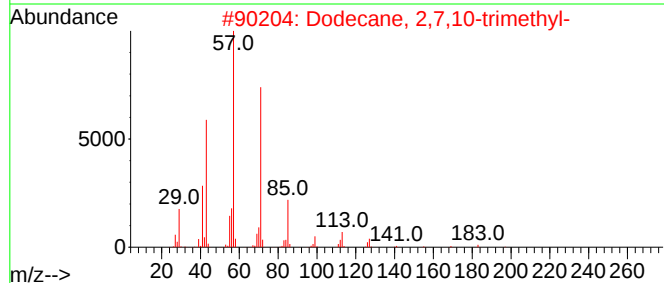
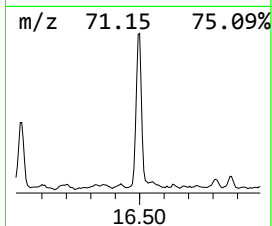
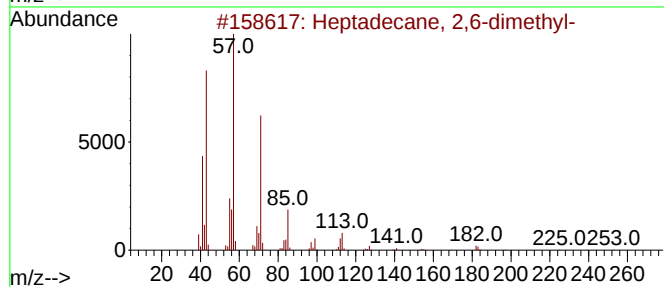
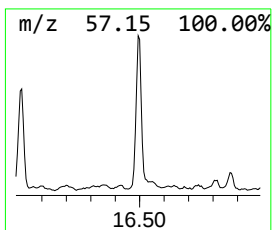
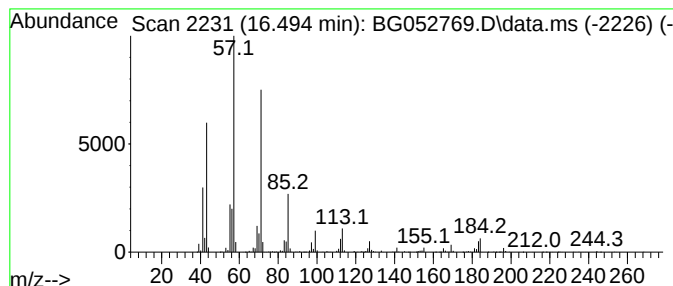
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Heptadecane, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.494	44.74 ng	1868070	Phenanthrene-d10	17.510	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	93
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	87
3	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	74
4	Tridecane	184	C13H28	000629-50-5	64
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

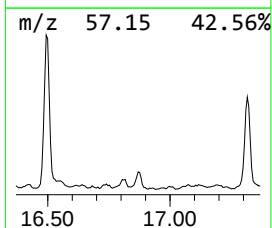
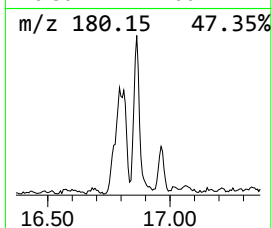
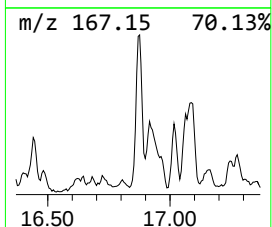
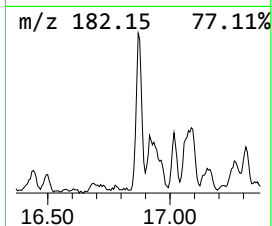
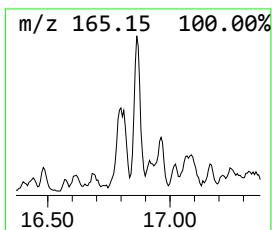
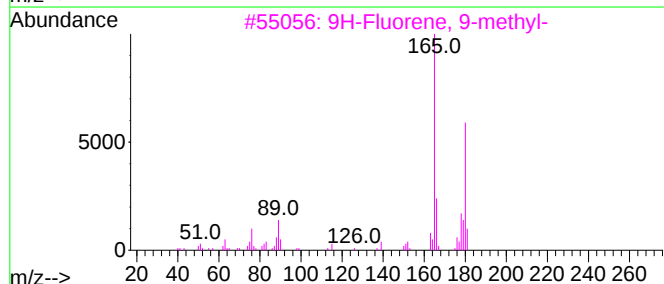
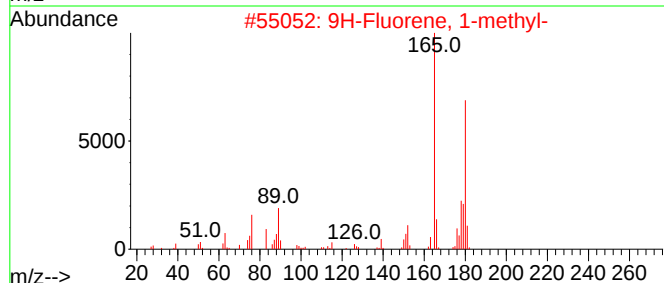
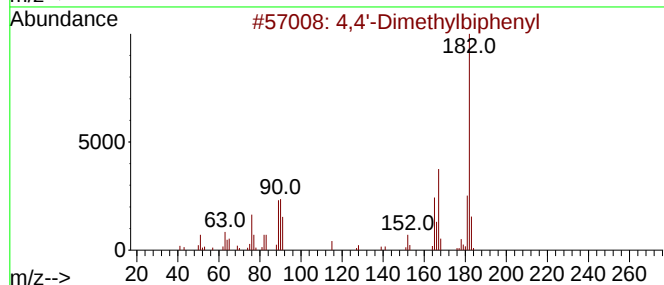
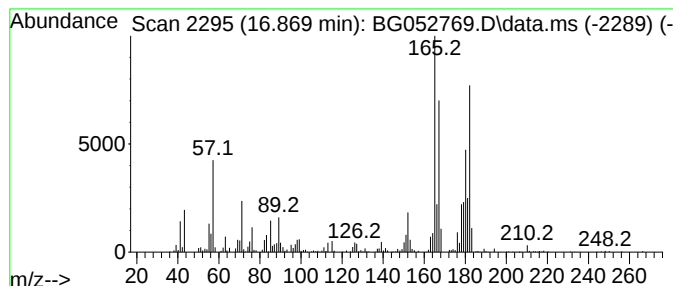
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 4,4'-Dimethylbiphenyl Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.869	17.22 ng	719016	Phenanthrene-d10		17.510	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4,4'-Dimethylbiphenyl		182	C14H14	000613-33-2	64
2	9H-Fluorene, 1-methyl-		180	C14H12	001730-37-6	49
3	9H-Fluorene, 9-methyl-		180	C14H12	002523-37-7	46
4	9H-Fluorene, 2-methyl-		180	C14H12	001430-97-3	46
5	3,3'-Dimethylbiphenyl		182	C14H14	000612-75-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

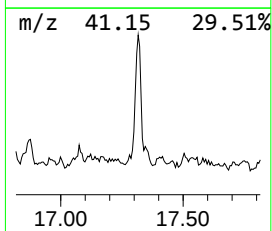
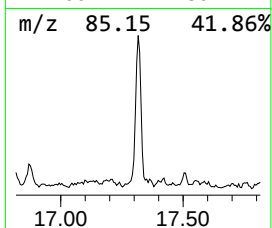
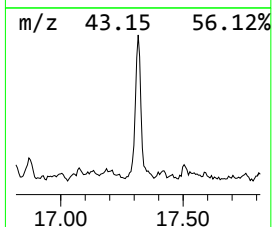
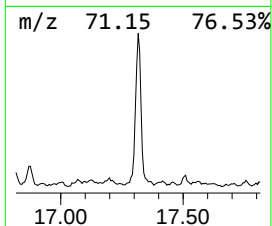
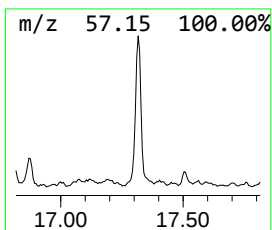
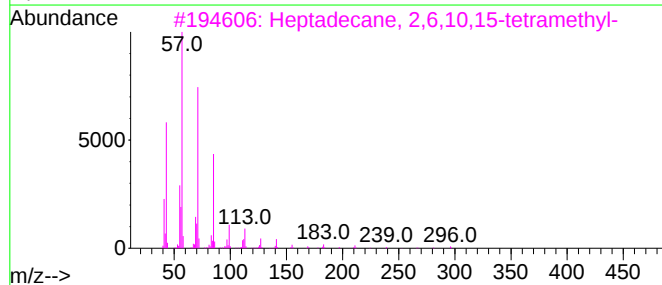
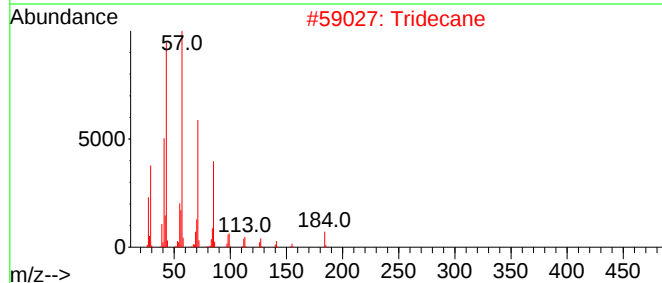
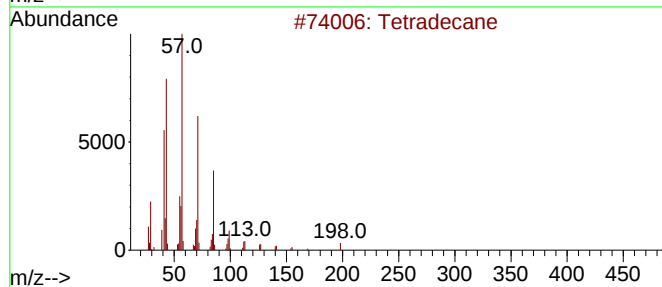
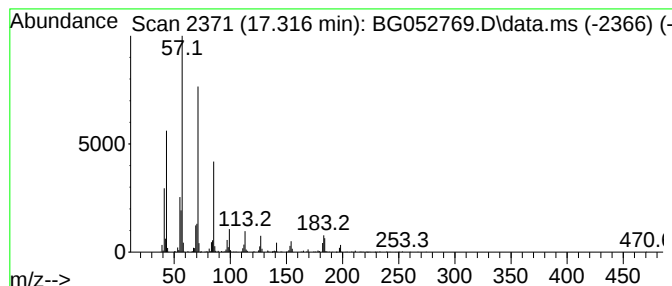
Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Tetradecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.316	34.03 ng	1420650	Phenanthrene-d10	17.510	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Tridecane	184	C13H28	000629-50-5	90
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
Data File : BG052769.D
Acq On : 21 Mar 2022 15:52
Operator : CG/JU
Sample : N1708-18 20X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PPSP-B10-25.5-26.0

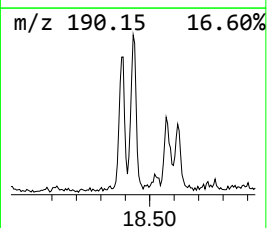
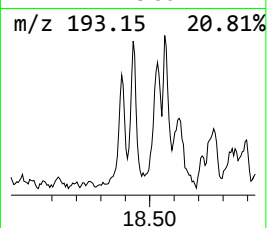
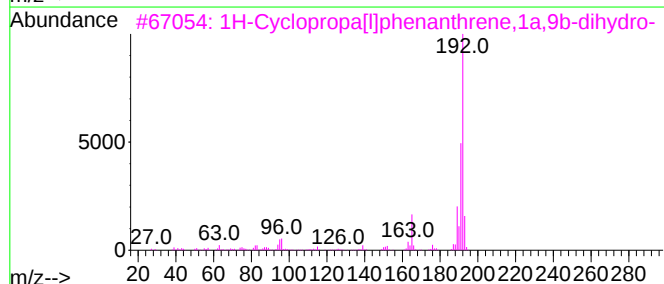
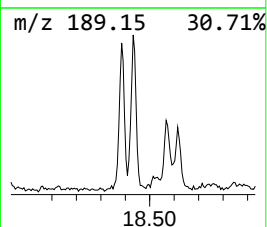
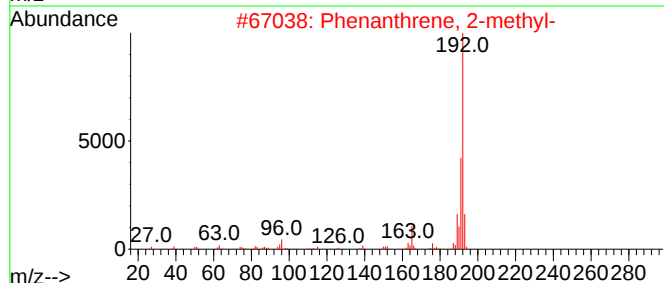
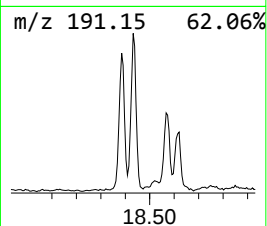
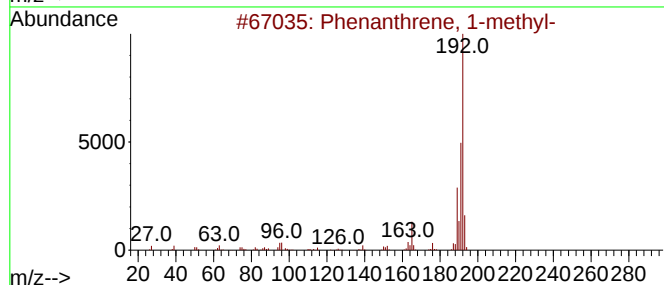
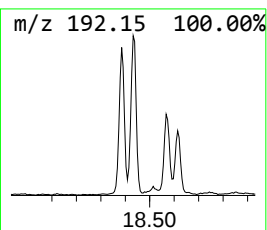
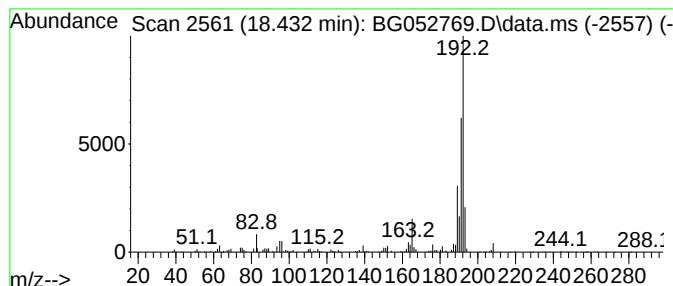
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Phenanthrene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.432	14.96 ng	624396	Phenanthrene-d10	17.510

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032122\
 Data File : BG052769.D
 Acq On : 21 Mar 2022 15:52
 Operator : CG/JU
 Sample : N1708-18 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PPSP-B10-25.5-26.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Naphthalene, 1,...	12.146	17.0	ng	996714	2	10.960	1173440	20.0
Nonane, 3,7-dim...	13.316	28.4	ng	960273	3	14.766	677364	20.0
Naphthalene, 1,...	13.715	15.2	ng	513996	3	14.766	677364	20.0
Naphthalene, 1,...	13.932	54.2	ng	1836990	3	14.766	677364	20.0
Naphthalene, 2,...	14.091	57.7	ng	1953540	3	14.766	677364	20.0
Pentadecane, 7-...	14.249	22.4	ng	759126	3	14.766	677364	20.0
Naphthalene, 2,...	14.326	30.1	ng	1017820	3	14.766	677364	20.0
Naphthalene, 2,...	14.972	29.8	ng	1010560	3	14.766	677364	20.0
Naphthalene, 1,...	15.160	27.8	ng	940414	3	14.766	677364	20.0
Naphthalene, 1,...	15.236	17.9	ng	606315	3	14.766	677364	20.0
unknown15.283	15.283	18.4	ng	624481	3	14.766	677364	20.0
Tridecane, 3-me...	16.018	31.6	ng	1071490	3	14.766	677364	20.0
Heptadecane, 2,...	16.494	44.7	ng	1868070	4	17.510	834990	20.0
4,4'-Dimethylbi...	16.869	17.2	ng	719016	4	17.510	834990	20.0
Tetradecane	17.316	34.0	ng	1420650	4	17.510	834990	20.0
Phenanthrene, 1...	18.432	15.0	ng	624396	4	17.510	834990	20.0