

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055482.D
 Acq On : 5 Nov 2022 23:11
 Operator : CG/JU
 Sample : N5459-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EO-03-110322

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-BG110122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055482.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.651	227	232	240	rBV2	8025	14800	1.70%	0.200%
2	5.274	330	338	346	rBV	72603	119156	13.69%	1.607%
3	5.762	414	421	434	rBV	256364	426438	48.98%	5.750%
4	7.384	690	697	713	rBV	246487	443010	50.88%	5.973%
5	8.230	835	841	850	rBV	135969	228295	26.22%	3.078%
6	9.405	1035	1041	1052	rBV	151887	276772	31.79%	3.732%
7	11.062	1316	1323	1329	rBV	191831	337611	38.78%	4.552%
8	11.114	1329	1332	1338	rVB	8209	12810	1.47%	0.173%
9	12.695	1597	1601	1609	rVB2	7469	10546	1.21%	0.142%
10	12.912	1632	1638	1646	rVB2	8397	16836	1.93%	0.227%
11	13.476	1726	1734	1745	rBV	458269	703599	80.82%	9.487%
12	14.029	1820	1828	1834	rBV6	8016	19153	2.20%	0.258%
13	14.170	1847	1852	1857	rBV2	14113	23840	2.74%	0.321%
14	14.223	1857	1861	1865	rVB2	9078	12194	1.40%	0.164%
15	14.411	1888	1893	1900	rVB5	7447	13434	1.54%	0.181%
16	14.581	1915	1922	1930	rBV5	8793	20741	2.38%	0.280%
17	14.851	1958	1968	1974	rBV	289886	470834	54.08%	6.348%
18	14.910	1974	1978	1984	rVB2	21246	32890	3.78%	0.443%
19	15.045	1995	2001	2002	rBV6	7300	10017	1.15%	0.135%
20	15.239	2026	2034	2040	rVB2	21740	43126	4.95%	0.581%
21	15.310	2042	2046	2051	rBV2	11534	16634	1.91%	0.224%
22	15.451	2064	2070	2075	rBV2	13009	26773	3.08%	0.361%
23	15.503	2075	2079	2085	rVB3	8145	16379	1.88%	0.221%
24	15.621	2093	2099	2102	rBV6	8656	17230	1.98%	0.232%
25	15.650	2102	2104	2109	rVB3	12709	14215	1.63%	0.192%
26	15.891	2135	2145	2149	rBV2	41051	79013	9.08%	1.065%
27	16.050	2168	2172	2176	rVB5	13314	19347	2.22%	0.261%
28	16.326	2214	2219	2228	rBV	317927	508985	58.46%	6.863%
29	16.526	2246	2253	2265	rBV7	19610	64676	7.43%	0.872%
30	17.090	2345	2349	2350	rBV4	8186	10565	1.21%	0.142%
31	17.295	2381	2384	2388	rBV3	14116	18151	2.08%	0.245%
32	17.348	2390	2393	2397	rVV2	18442	23345	2.68%	0.315%
33	17.401	2399	2402	2407	rVB	24295	36227	4.16%	0.488%
34	17.583	2427	2433	2437	rBV	336215	528703	60.73%	7.128%
35	17.624	2437	2440	2447	rVB	292394	412367	47.37%	5.560%
36	17.713	2451	2455	2462	rVV	61321	100122	11.50%	1.350%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

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-BG110122.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	17.983	2498	2501	2506	rVB	31488	42500	4.88%	0.573%
38	18.030	2506	2509	2516	rVB3	17473	24668	2.83%	0.333%
39	18.159	2528	2531	2537	rBV	23115	32341	3.71%	0.436%
40	18.453	2575	2581	2585	rBV2	52785	114172	13.11%	1.539%
41	18.500	2586	2589	2593	rVB	58764	81812	9.40%	1.103%
42	18.647	2609	2614	2618	rBV2	53925	110785	12.72%	1.494%
43	19.616	2775	2779	2785	rVB	391157	524099	60.20%	7.066%
44	19.980	2837	2841	2846	rVB	346323	487008	55.94%	6.566%
45	20.157	2867	2871	2876	rVB	694540	870614	100.00%	11.738%

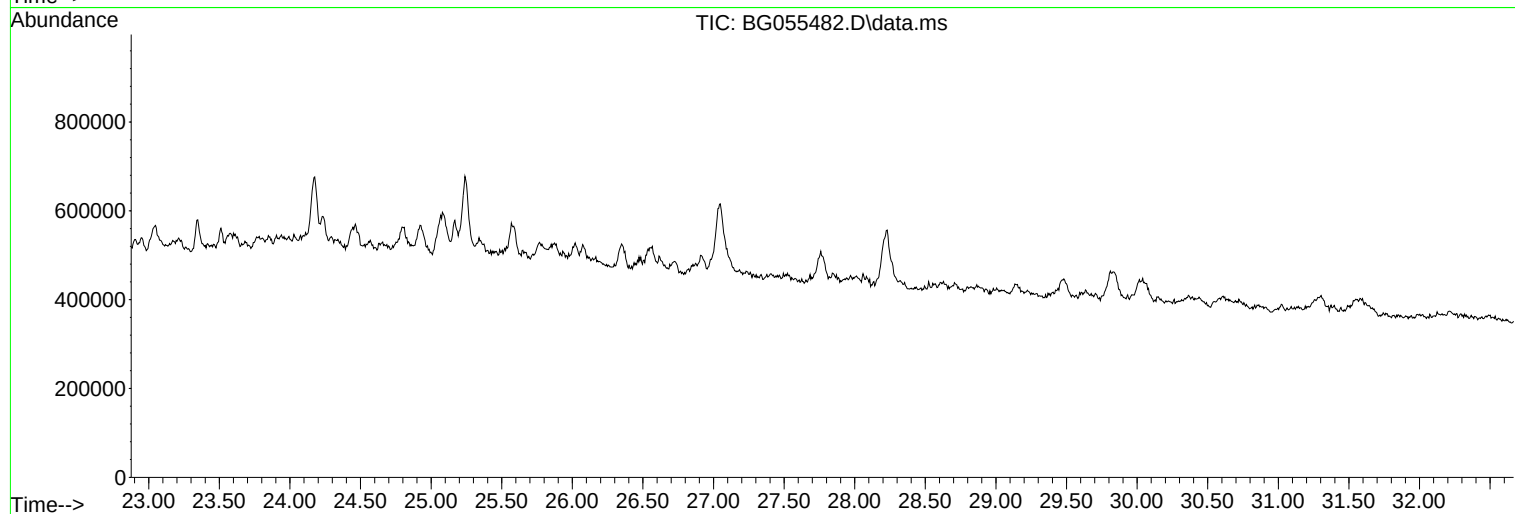
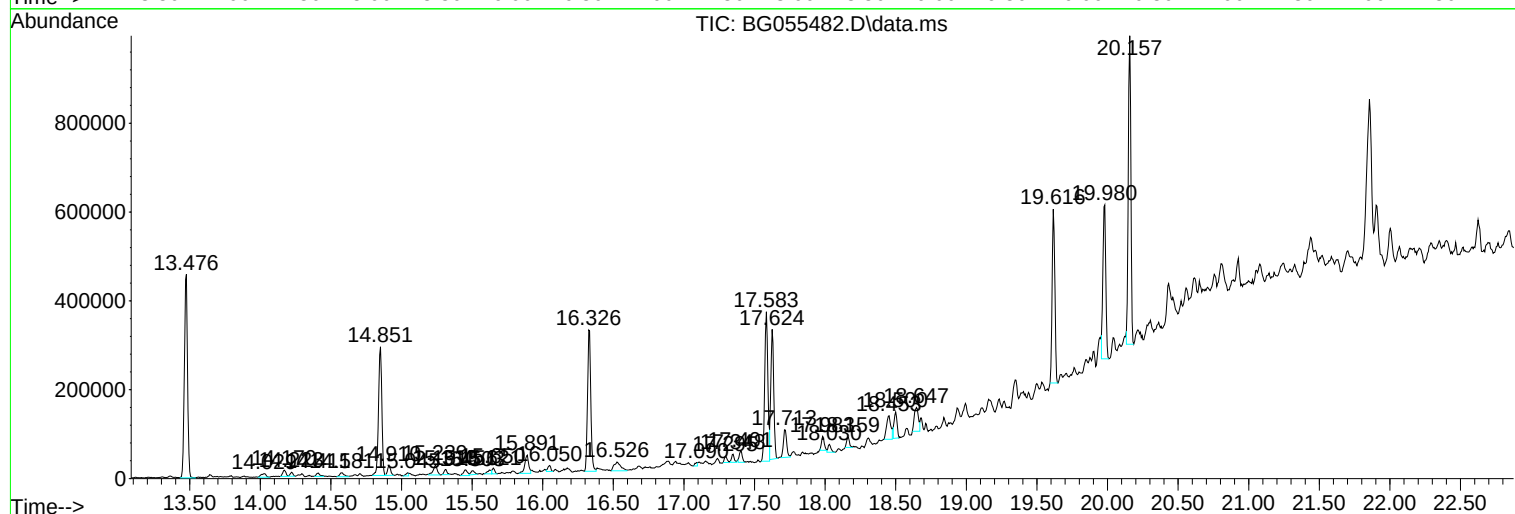
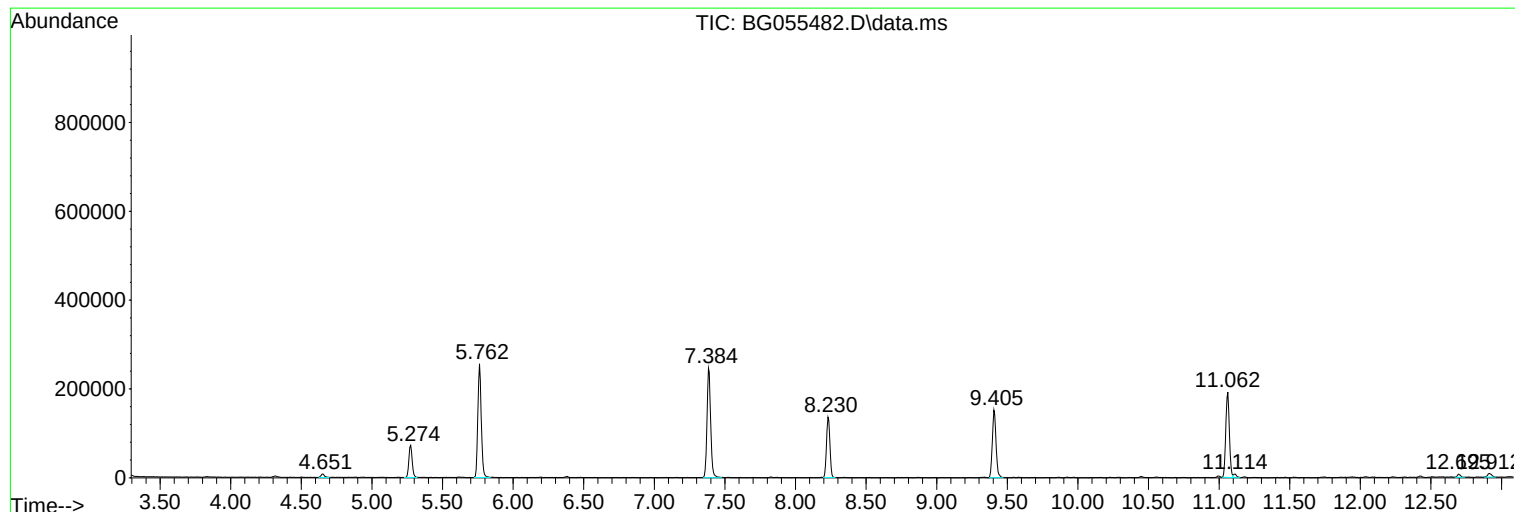
Sum of corrected areas: 7416833

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.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\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

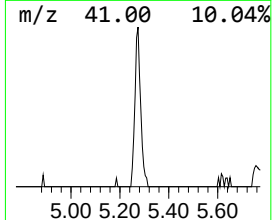
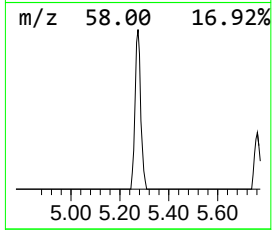
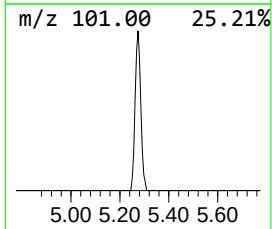
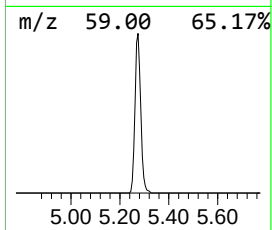
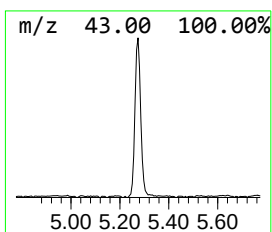
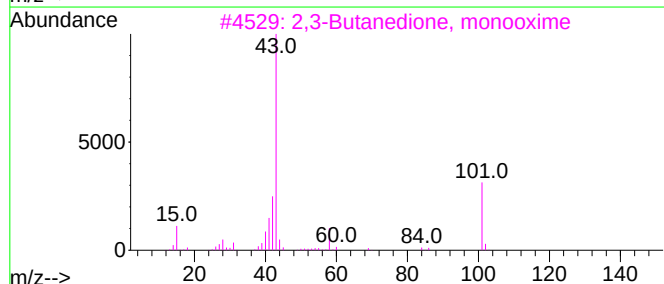
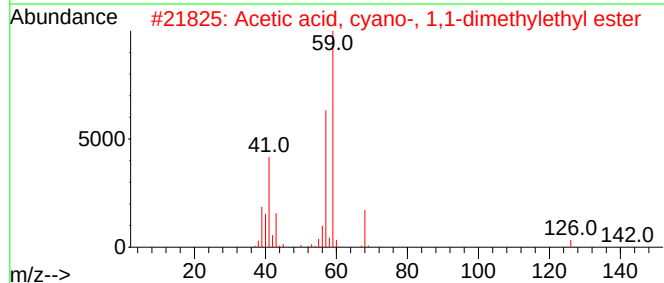
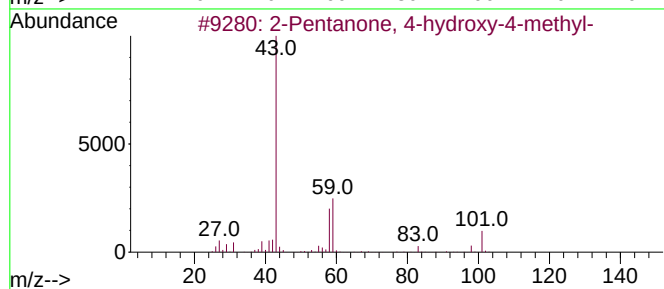
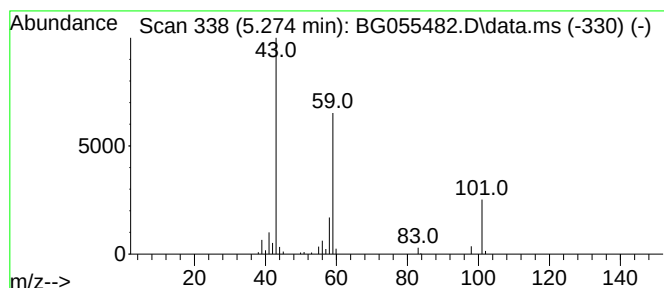
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.274	10.44 ng	119156	1,4-Dichlorobenzene-d4	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

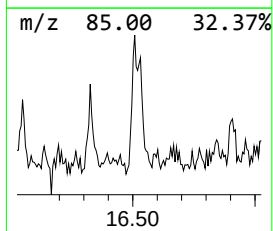
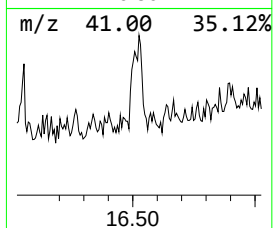
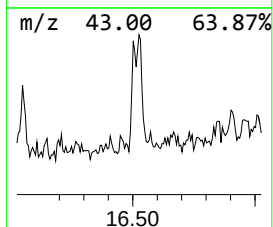
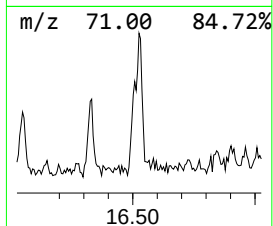
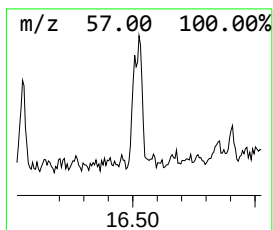
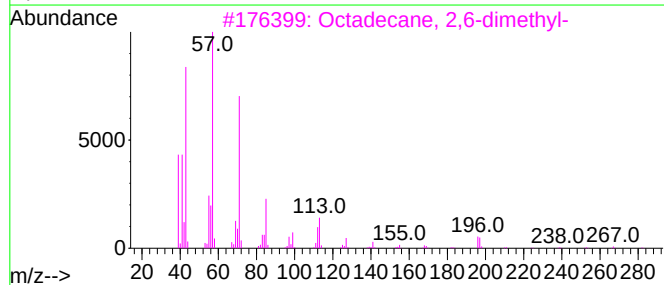
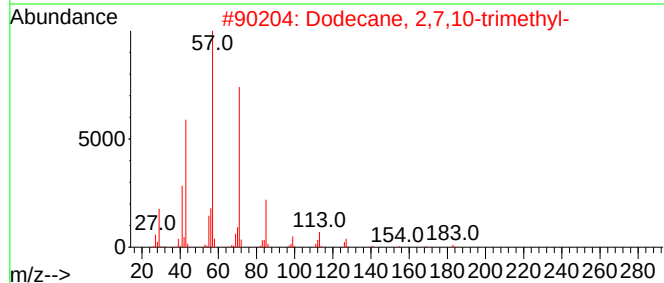
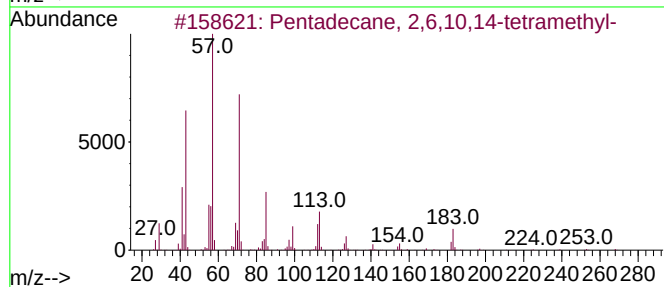
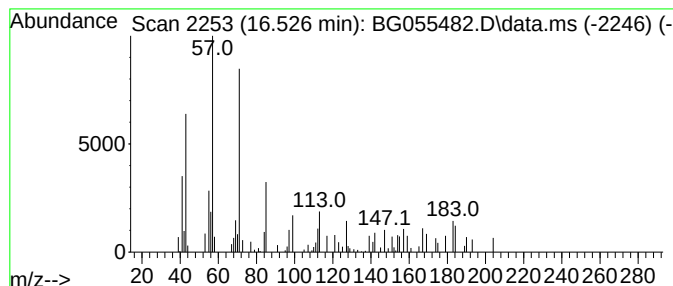
Instrument :
BNA_G
ClientSampleId :
EO-03-110322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Pentadecane, 2,6,10,14-tetr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.526	2.45 ng	64676	Phenanthrene-d10		17.583	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	64
2	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	59
3	Octadecane, 2,6-dimethyl-		282	C20H42	075163-97-2	53
4	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	53
5	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

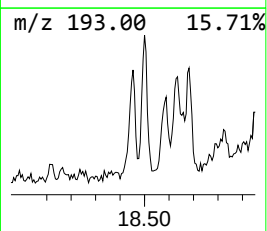
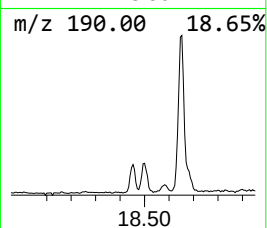
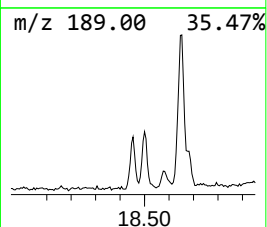
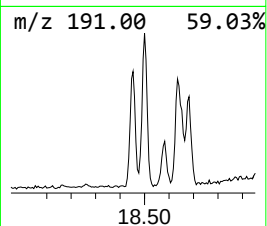
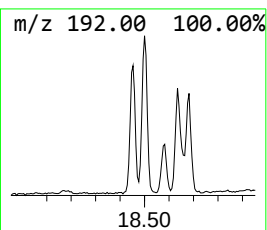
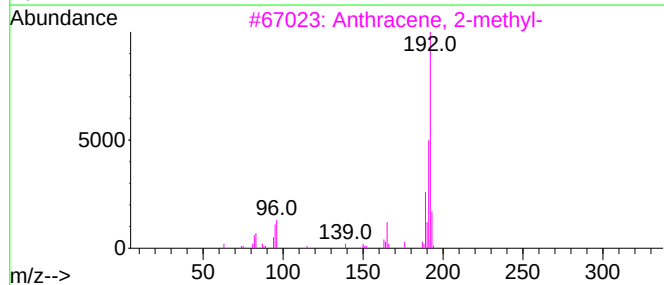
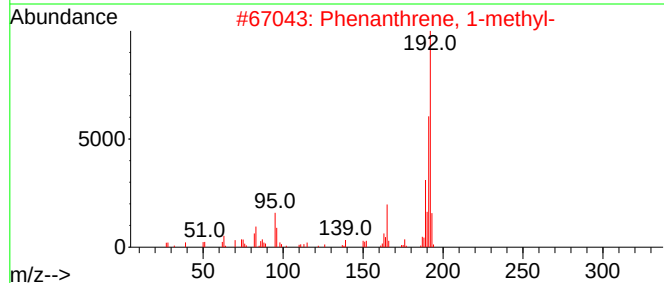
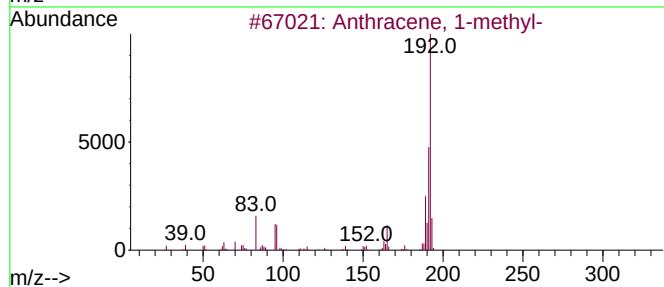
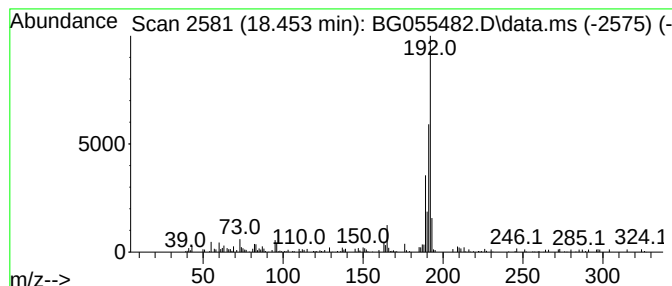
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Anthracene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.453	4.32 ng	114172	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

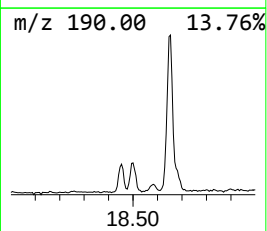
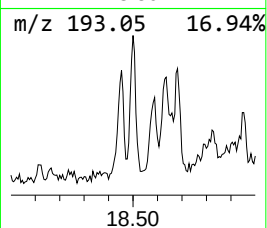
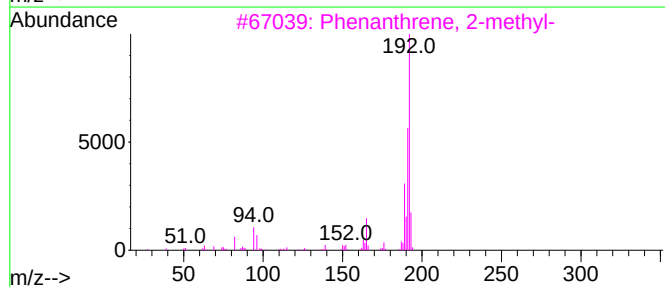
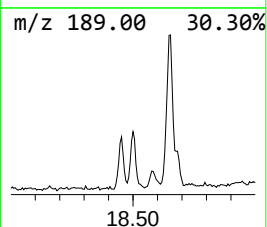
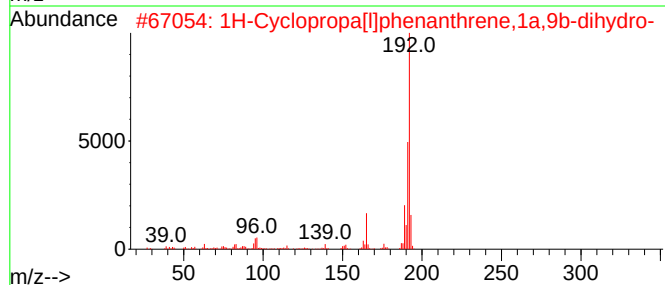
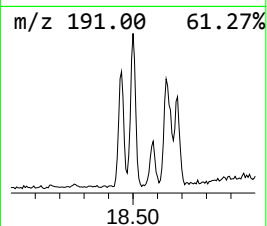
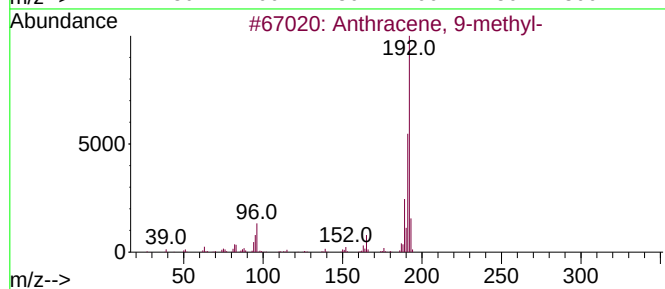
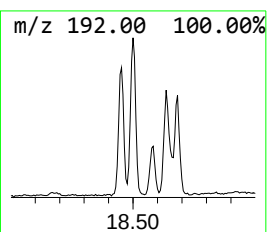
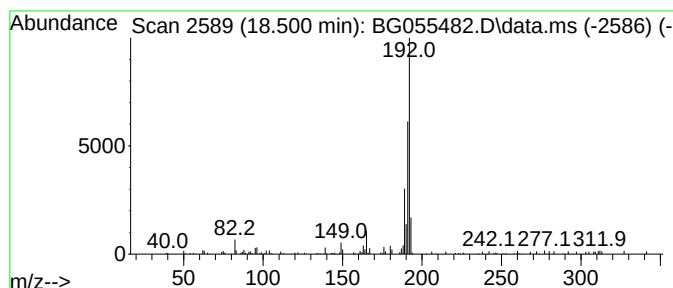
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Anthracene, 9-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.500	3.09 ng	81812	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	87
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

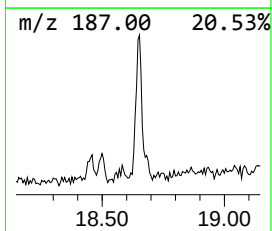
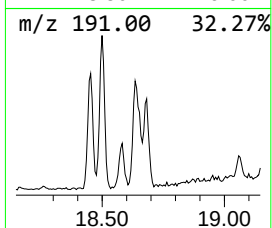
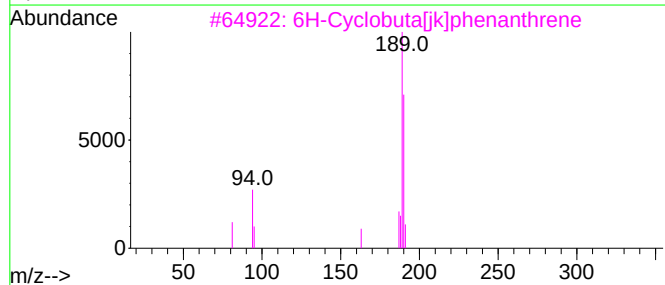
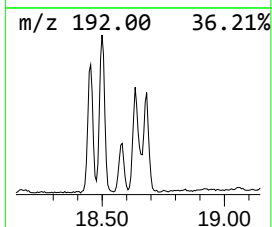
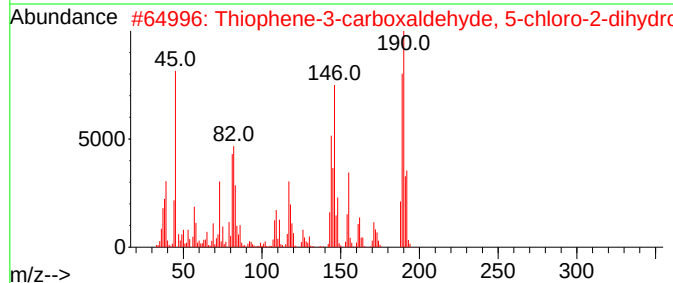
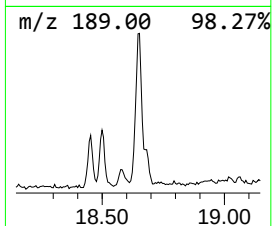
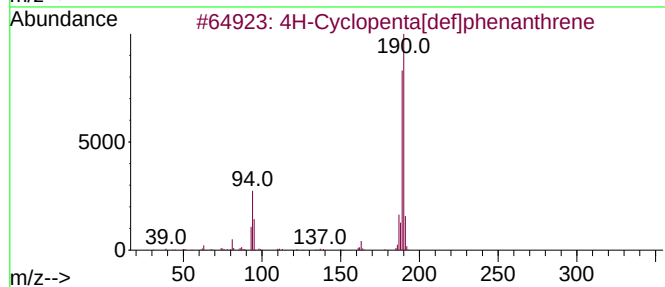
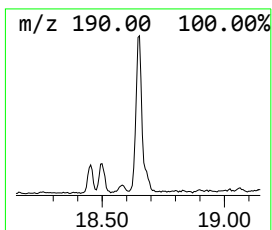
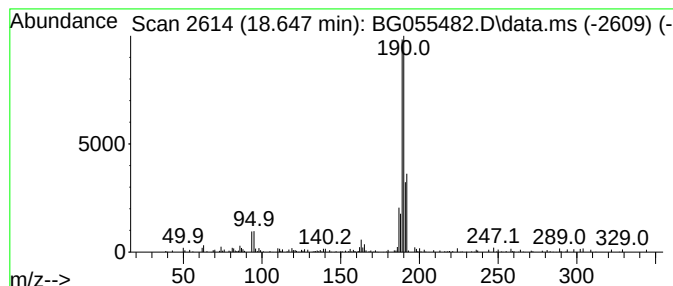
Instrument :
BNA_G
ClientSampleId :
EO-03-110322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.647	4.19 ng	110785	Phenanthrene-d10		17.583	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	74
2	Thiophene-3-carboxaldehyde, 5-ch...		190	C5H4BClO3S	036155-87-0	72
3	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	72
4	2,3,5,6-Tetramethylterephthalald...		190	C12H14O2	007072-01-7	50
5	Phenol, 4-(2-thienylmethyl)-		190	C11H10O5	091680-55-6	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
Data File : BG055482.D
Acq On : 5 Nov 2022 23:11
Operator : CG/JU
Sample : N5459-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.274	10.4	ng	119156	1	8.230	228295	20.0
Pentadecane, 2,...	16.526	2.5	ng	64676	4	17.583	528703	20.0
Anthracene, 1-m...	18.453	4.3	ng	114172	4	17.583	528703	20.0
Anthracene, 9-m...	18.500	3.1	ng	81812	4	17.583	528703	20.0
4H-Cyclopenta[d...	18.647	4.2	ng	110785	4	17.583	528703	20.0