

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
 Data File : BG054422.D
 Acq On : 27 Jul 2022 19:10
 Operator : CG/JU
 Sample : N3892-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BIN-4

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-BG071522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG054422.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.032	3	7	15	rVB4	9789	19392	1.79%	0.245%
2	3.108	15	20	30	rVB	67460	117579	10.88%	1.487%
3	5.564	431	438	447	rBV	183838	320144	29.61%	4.048%
4	7.174	704	712	725	rBV	197568	356998	33.02%	4.514%
5	7.991	844	851	858	rBV	60577	106960	9.89%	1.352%
6	9.154	1041	1049	1057	rBV	127632	233682	21.61%	2.955%
7	10.799	1318	1329	1334	rBV	79931	148039	13.69%	1.872%
8	10.846	1334	1337	1346	rVB2	11133	17679	1.64%	0.224%
9	12.450	1604	1610	1618	rVB	8252	14633	1.35%	0.185%
10	13.244	1738	1745	1754	rBV	414635	671800	62.14%	8.494%
11	13.943	1858	1864	1871	rBV3	6288	12012	1.11%	0.152%
12	14.348	1927	1933	1945	rBV	13107	27829	2.57%	0.352%
13	14.513	1955	1961	1968	rVB2	8020	11584	1.07%	0.146%
14	14.624	1968	1980	1986	rBV	143945	235727	21.80%	2.980%
15	14.689	1986	1991	1998	rVB	11342	19625	1.82%	0.248%
16	15.024	2043	2048	2055	rBV	23302	36864	3.41%	0.466%
17	15.465	2118	2123	2126	rVB2	9446	13407	1.24%	0.170%
18	15.670	2152	2158	2170	rBV	32656	58345	5.40%	0.738%
19	15.964	2202	2208	2215	rVB	9843	17340	1.60%	0.219%
20	16.117	2226	2234	2243	rBV2	379364	615999	56.98%	7.788%
21	16.322	2265	2269	2270	rBV	15988	17144	1.59%	0.217%
22	16.346	2271	2273	2278	rVB3	19238	22614	2.09%	0.286%
23	16.675	2320	2329	2334	rBV5	11245	28396	2.63%	0.359%
24	16.722	2334	2337	2342	rVB3	10213	15832	1.46%	0.200%
25	16.822	2349	2354	2360	rVB7	7484	14443	1.34%	0.183%
26	16.892	2360	2366	2371	rBV7	13738	27343	2.53%	0.346%
27	17.110	2398	2403	2408	rBV3	21639	34039	3.15%	0.430%
28	17.186	2408	2416	2422	rVB3	20261	52952	4.90%	0.669%
29	17.368	2441	2447	2451	rBV	180878	299939	27.74%	3.792%
30	17.409	2451	2454	2462	rVV	234090	363863	33.65%	4.600%
31	17.503	2464	2470	2475	rVV	71140	110290	10.20%	1.394%
32	17.562	2475	2480	2485	rVV4	9642	19664	1.82%	0.249%
33	17.644	2491	2494	2499	rVB6	7731	11878	1.10%	0.150%
34	17.779	2513	2517	2521	rBV3	13087	17314	1.60%	0.219%
35	17.850	2526	2529	2533	rBV	21913	30309	2.80%	0.383%
36	18.249	2592	2597	2601	rBV2	51372	90408	8.36%	1.143%

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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-BG071522.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	18.291	2601	2604	2613	rVB2	49164	84999	7.86%	1.075%
38	18.373	2614	2618	2623	rBV2	18936	33140	3.07%	0.419%
39	18.443	2624	2630	2634	rBV3	76485	141375	13.08%	1.787%
40	18.543	2644	2647	2651	rBV2	28754	34117	3.16%	0.431%
41	18.743	2677	2681	2684	rVB	27604	36584	3.38%	0.463%
42	18.790	2686	2689	2695	rVB3	21222	33316	3.08%	0.421%
43	19.166	2748	2753	2757	rBV3	45605	89638	8.29%	1.133%
44	19.425	2792	2797	2802	rVB	409313	586565	54.25%	7.416%
45	19.754	2849	2853	2855	rBV2	78739	123599	11.43%	1.563%
46	19.789	2855	2859	2867	rVB	366146	552980	51.15%	6.992%
47	19.977	2886	2891	2896	rVB	777837	1081169	100.00%	13.670%
48	20.277	2939	2942	2946	rVB	107642	150289	13.90%	1.900%
49	21.634	3166	3173	3178	rBV2	282103	749441	69.32%	9.475%

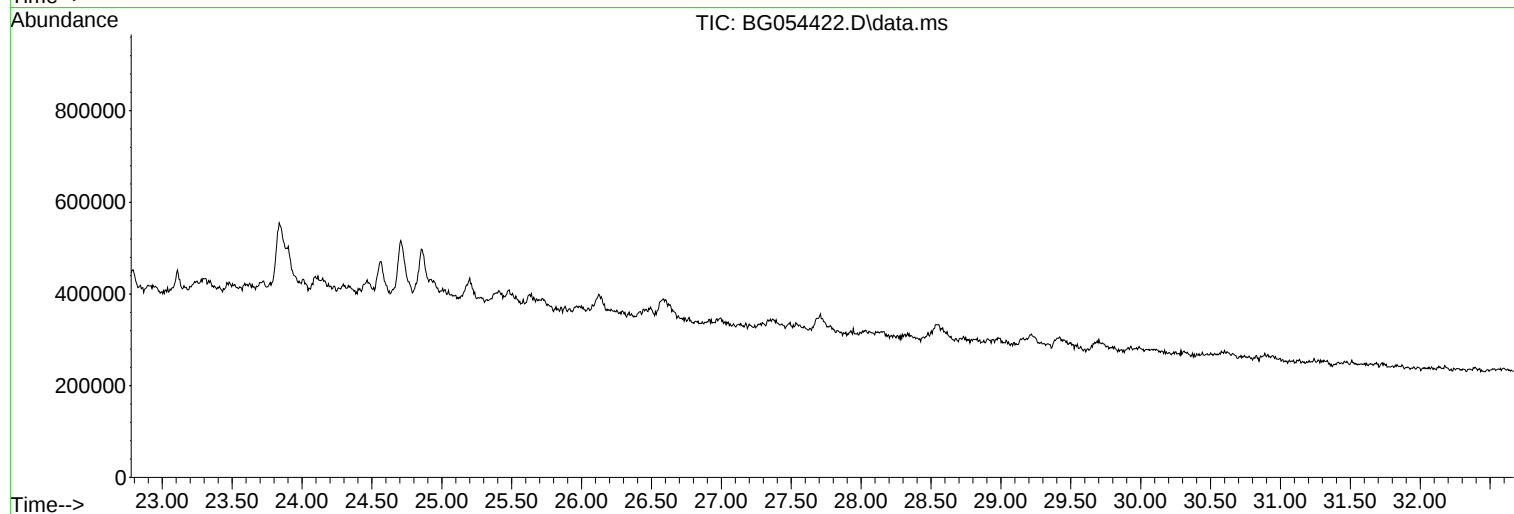
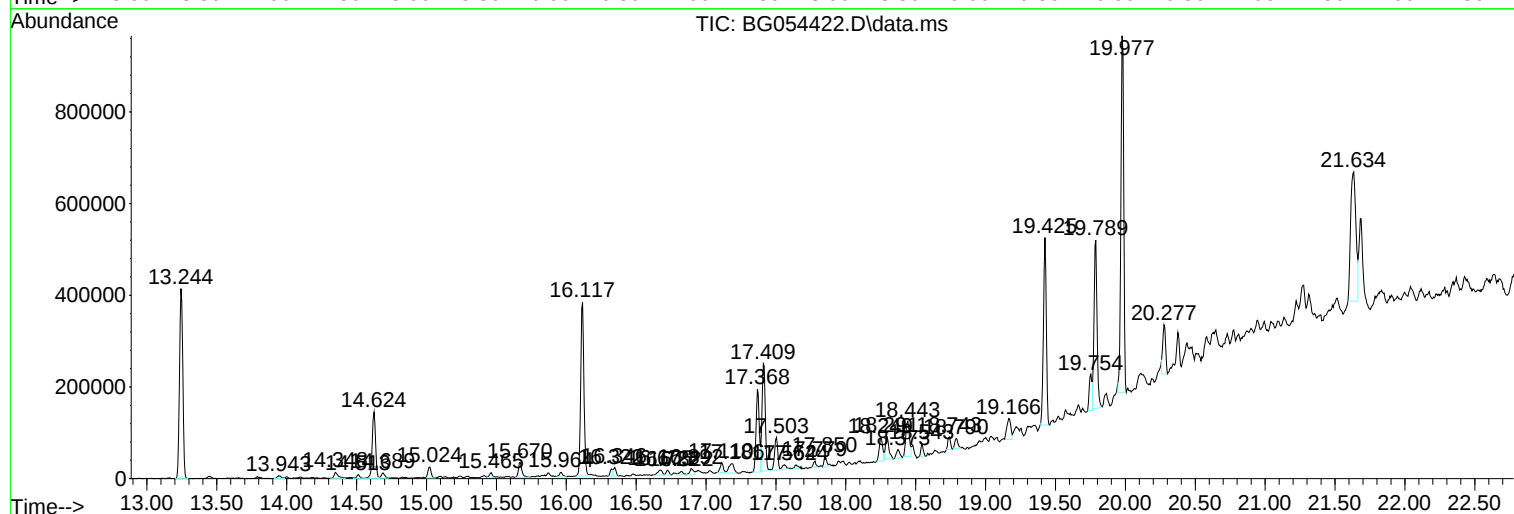
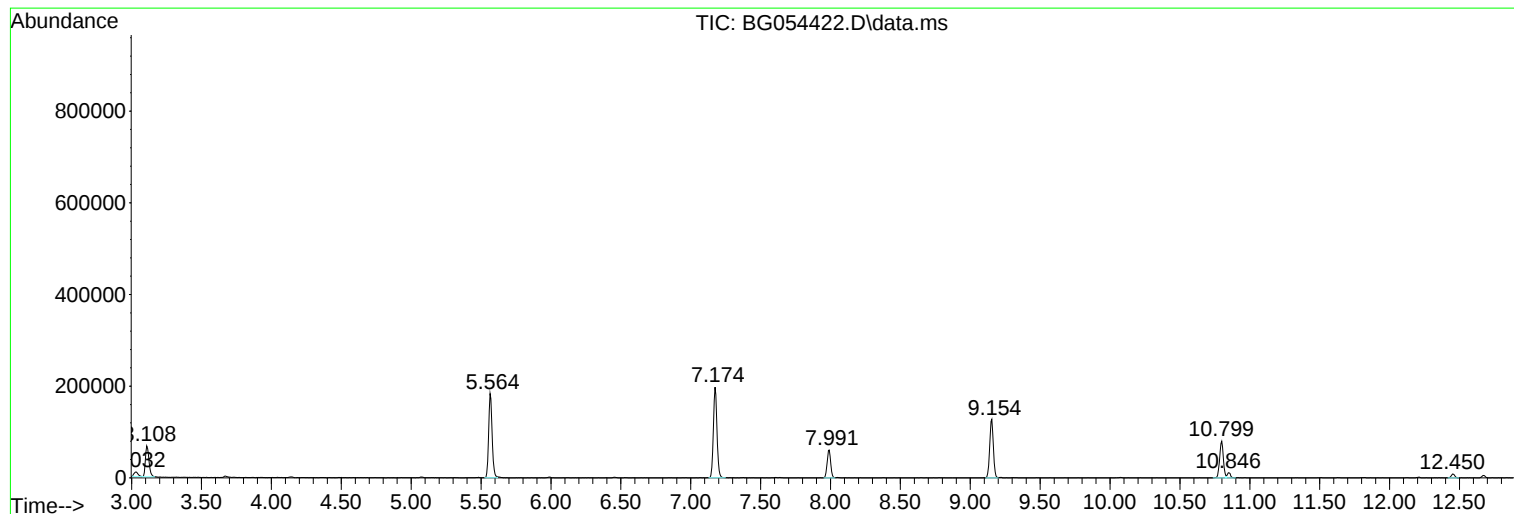
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ClientSampleId :
BIN-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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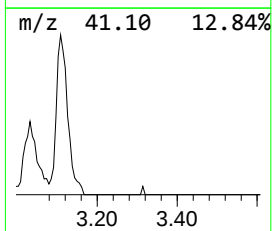
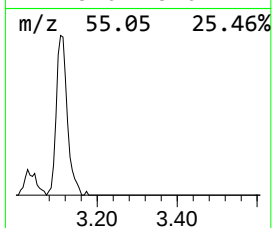
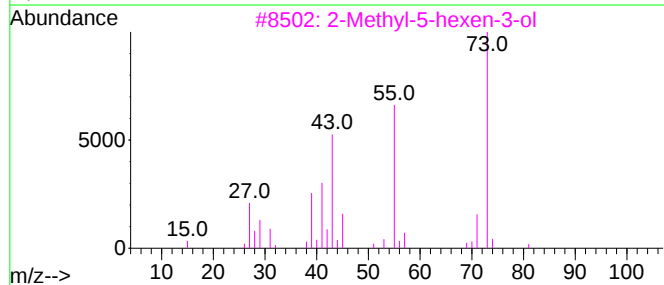
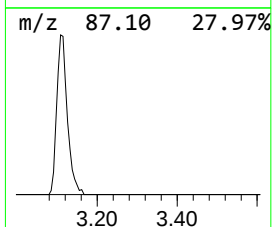
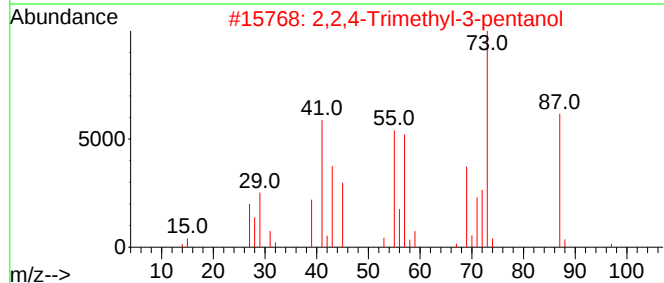
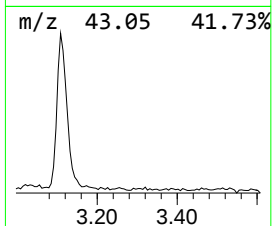
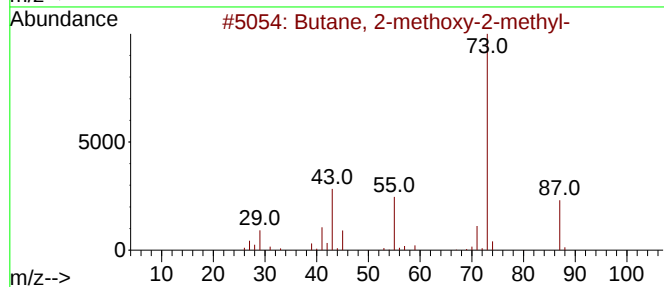
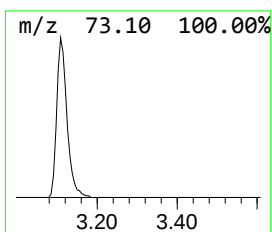
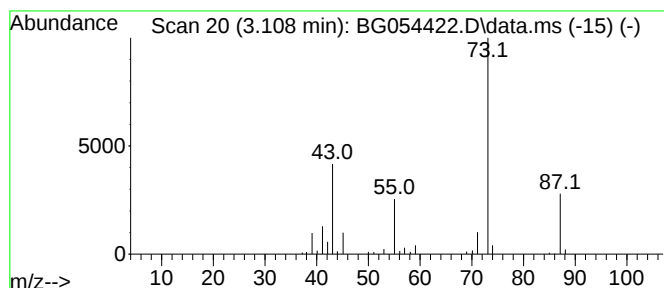
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.108	21.99 ng	117579	1,4-Dichlorobenzene-d4	7.991

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	50
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	16
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9



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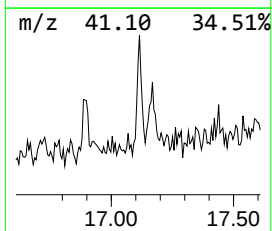
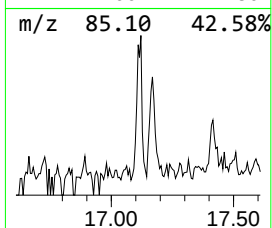
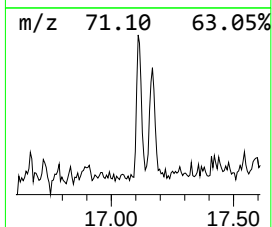
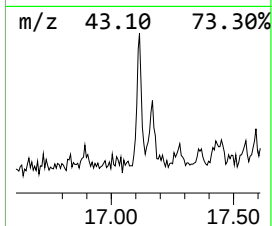
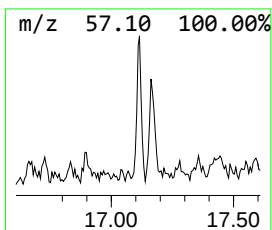
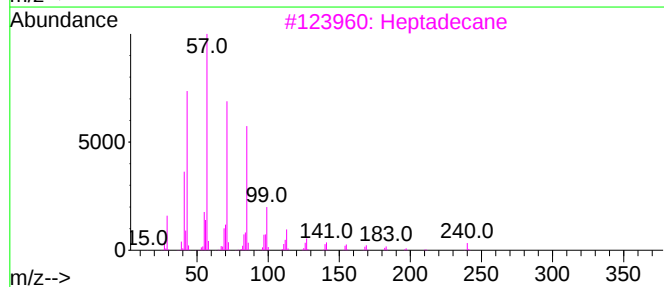
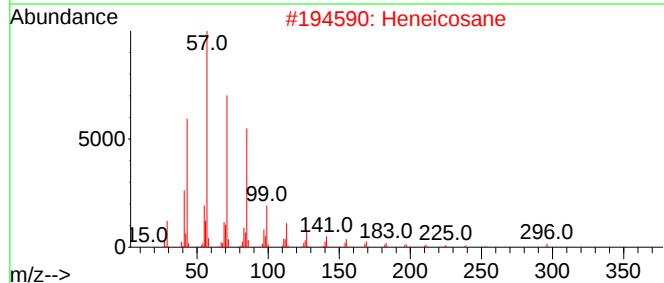
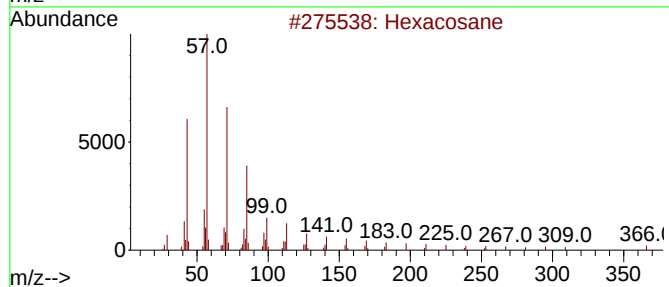
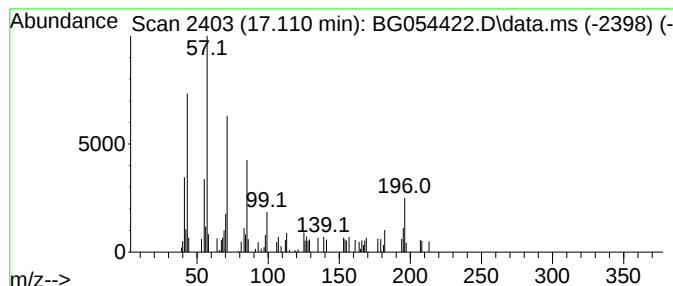
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.110	2.27 ng	34039	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	58
2			Heneicosane	296	C21H44	000629-94-7	58
3			Heptadecane	240	C17H36	000629-78-7	58
4			Octadecane	254	C18H38	000593-45-3	52
5			Pentacosane	352	C25H52	000629-99-2	52



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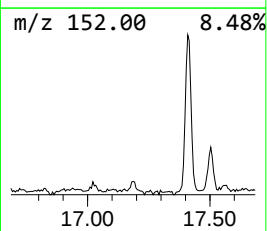
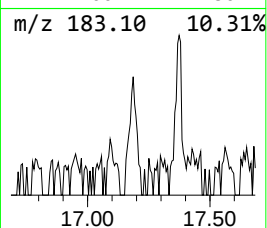
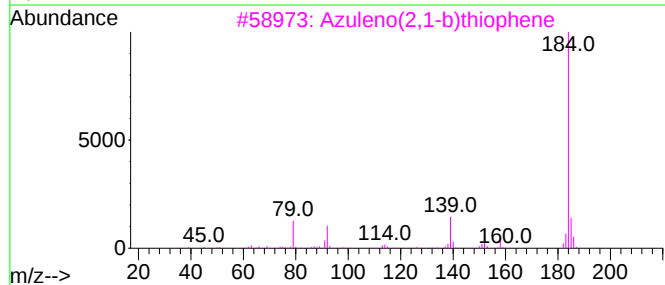
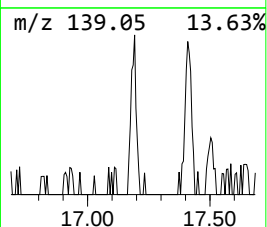
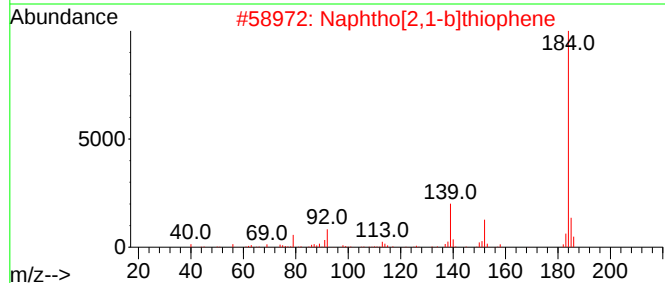
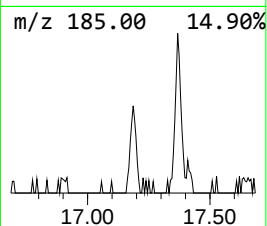
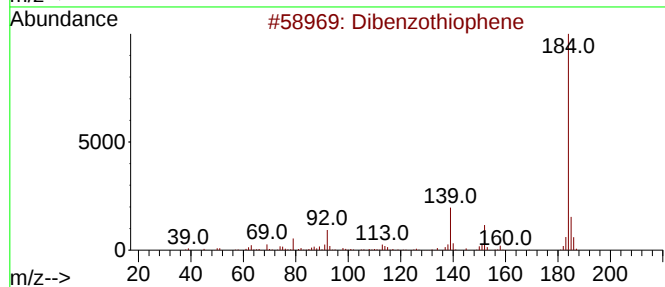
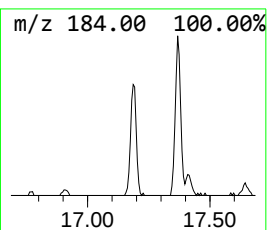
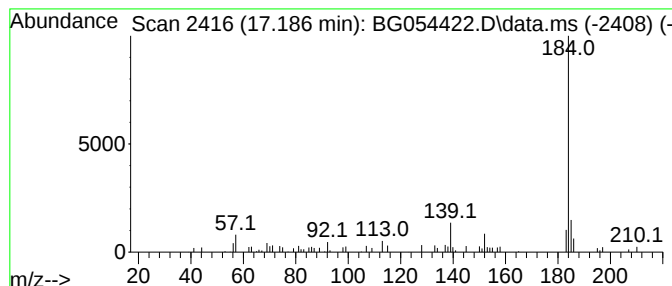
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Dibenzothiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.186	3.53 ng	52952	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	91
2			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	91
3			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	83
4			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	80
5			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	72



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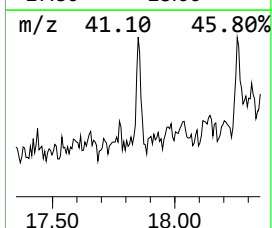
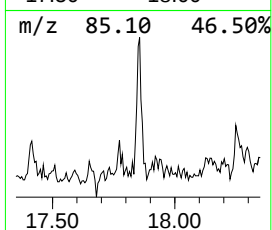
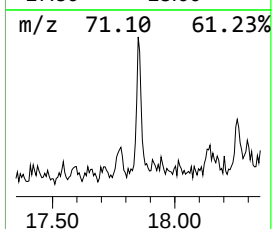
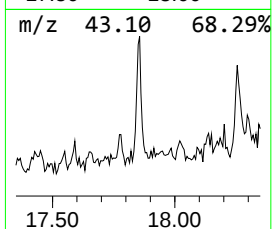
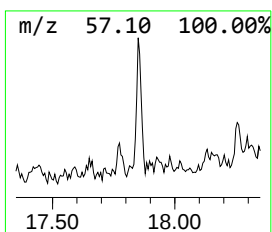
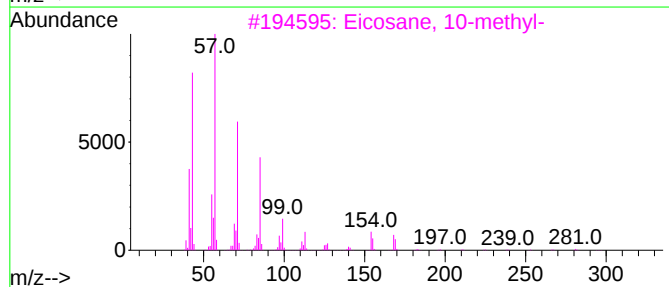
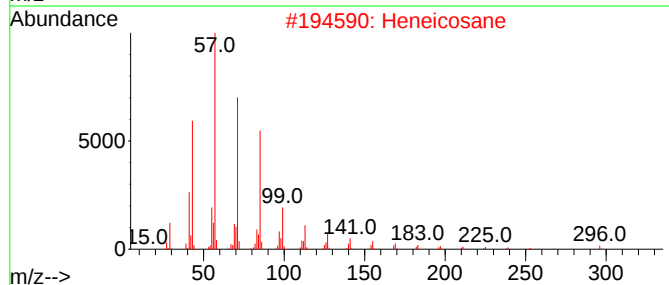
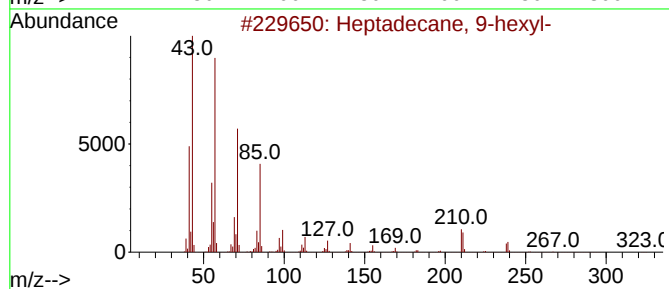
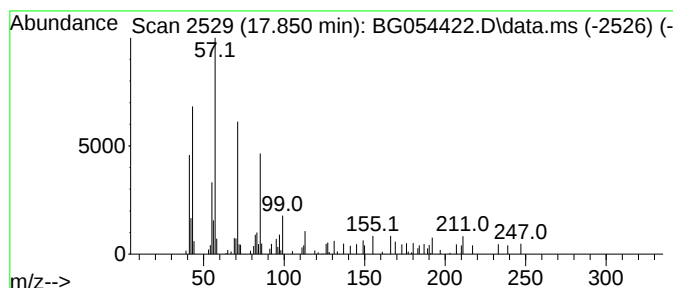
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Heptadecane, 9-hexyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	2.02 ng	30309	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	74
2			Heneicosane	296	C21H44	000629-94-7	74
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
4			Octacosane	394	C28H58	000630-02-4	72
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	72



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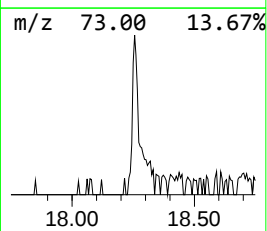
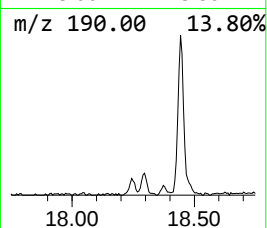
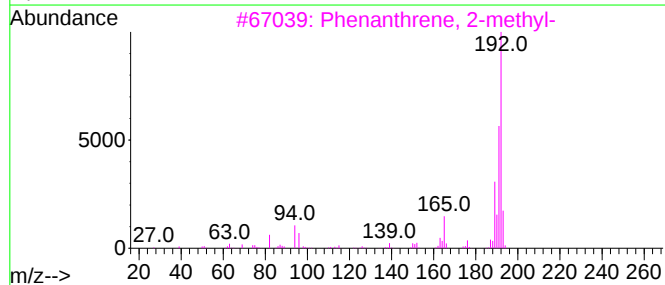
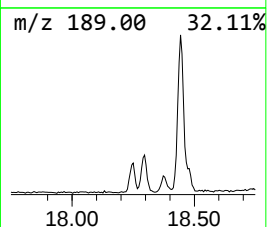
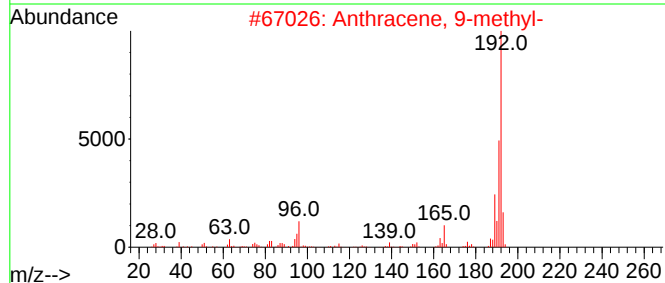
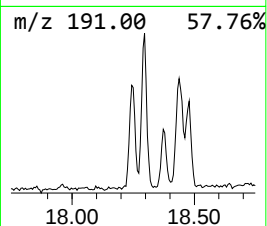
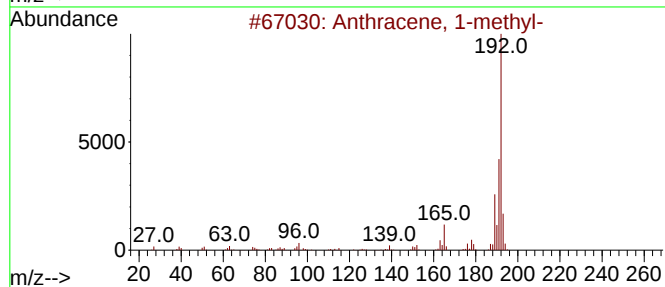
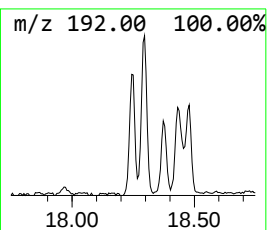
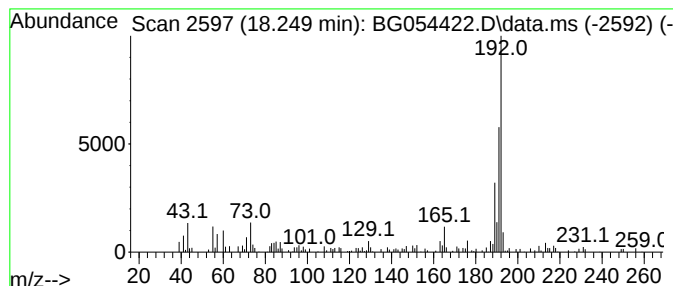
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.250	6.03 ng	90408	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	86
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	86
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	86
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
Operator : CG/JU
Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BIN-4

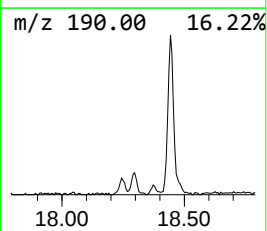
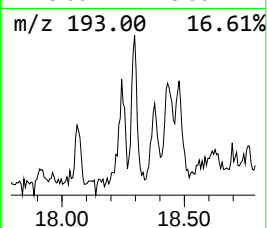
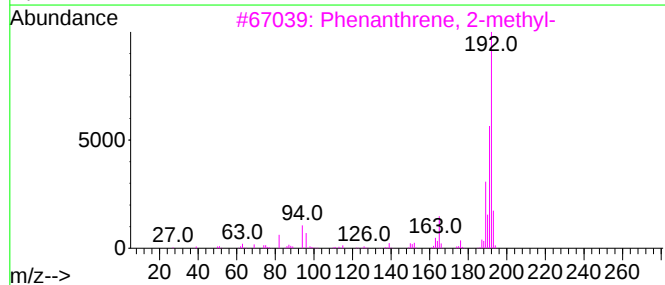
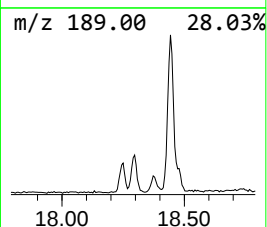
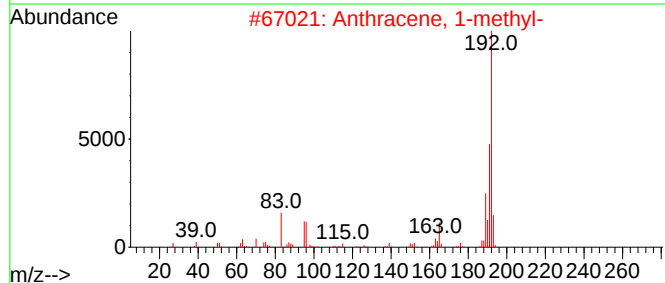
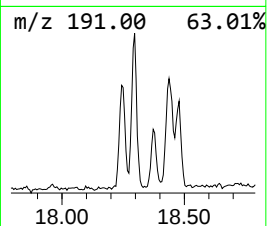
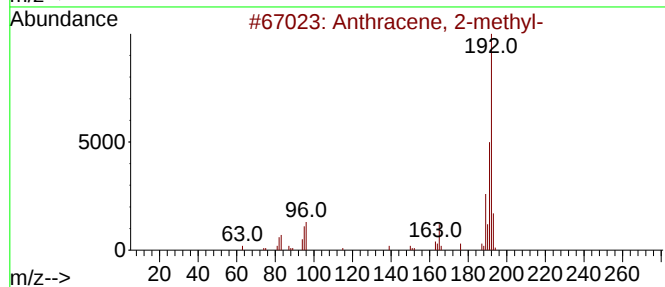
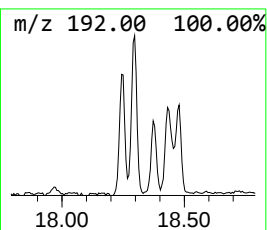
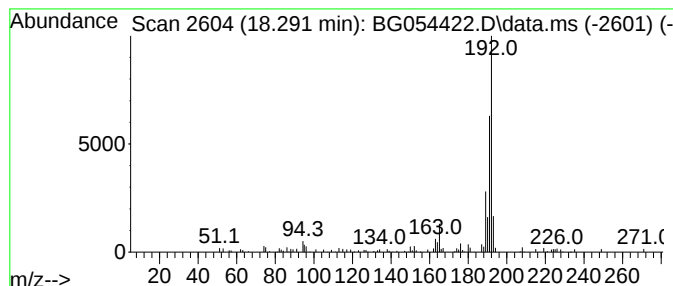
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.291	5.67 ng	84999	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
Operator : CG/JU
Sample : N3892-01
Misc :
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Instrument :
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ClientSampleId :
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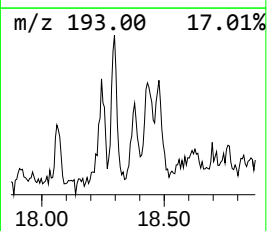
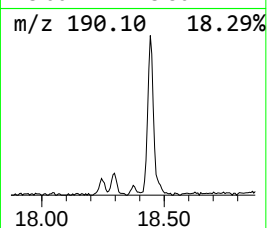
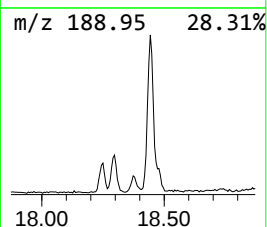
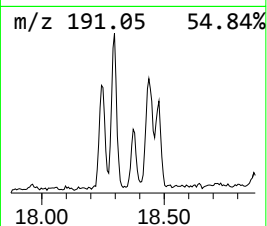
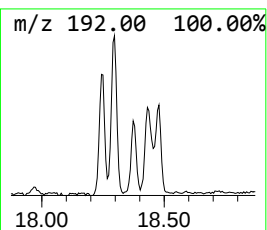
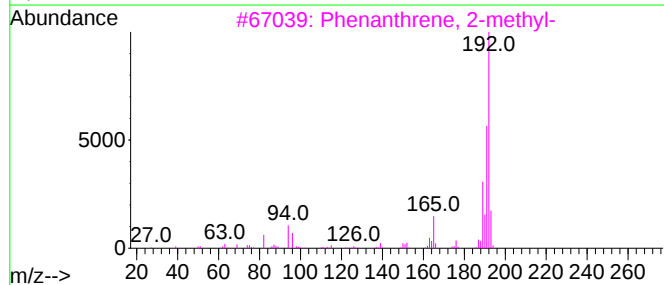
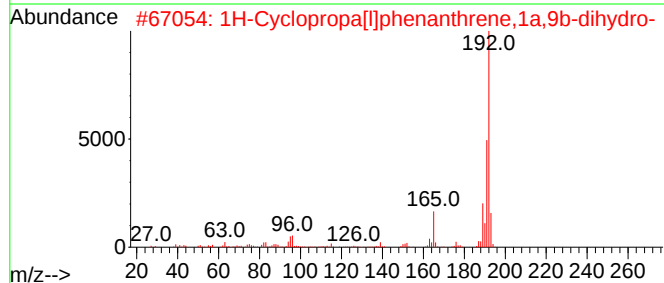
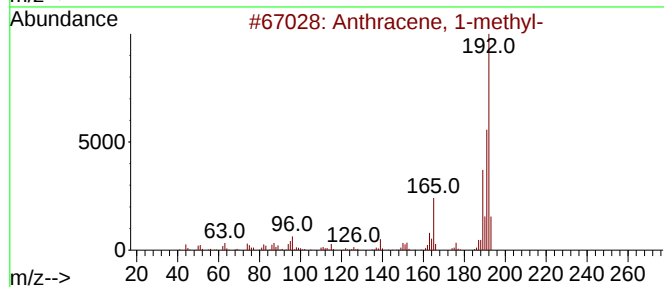
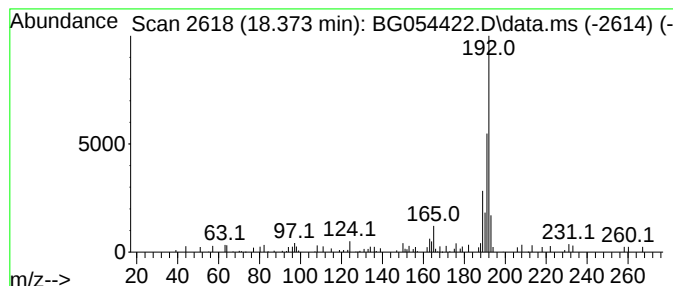
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.373	2.21 ng	33140	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
2			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	87
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
4			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	87
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
Operator : CG/JU
Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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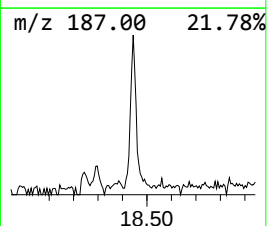
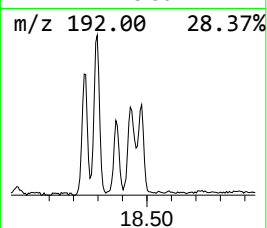
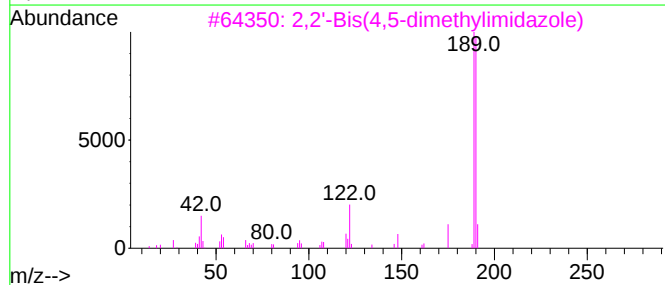
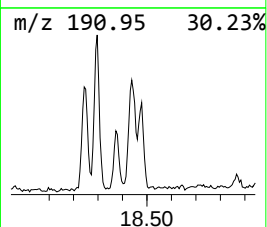
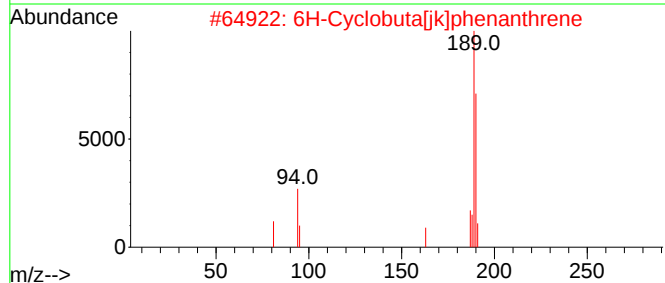
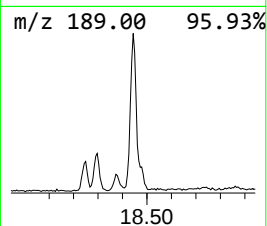
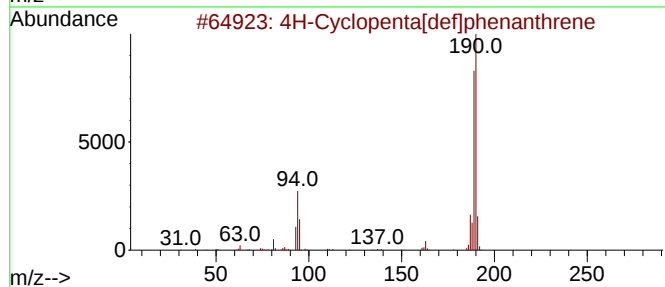
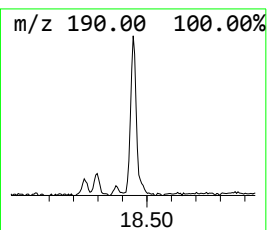
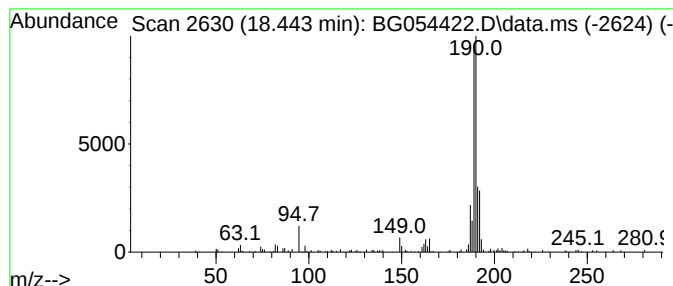
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.443	9.43 ng	141375	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	36
4			4-(4-Chlorophenyl)pyridine	189	C11H8ClN	005957-96-0	35
5			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
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Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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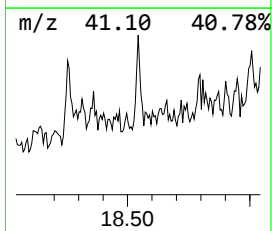
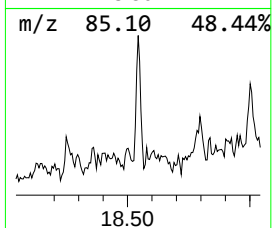
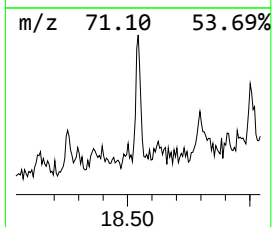
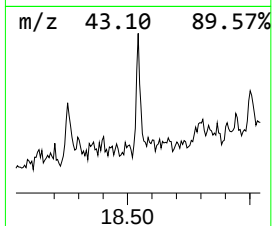
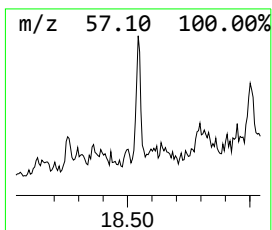
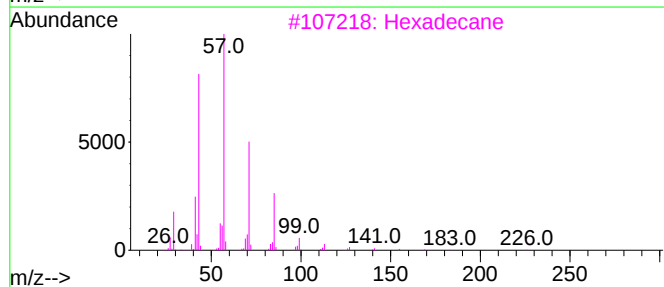
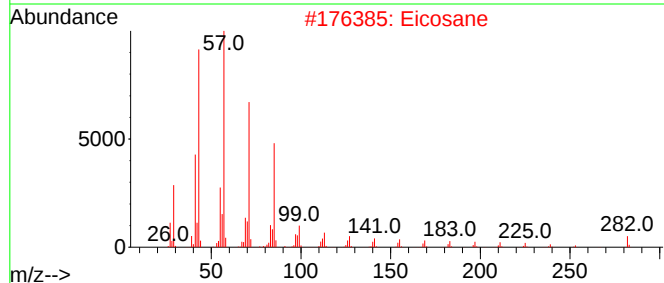
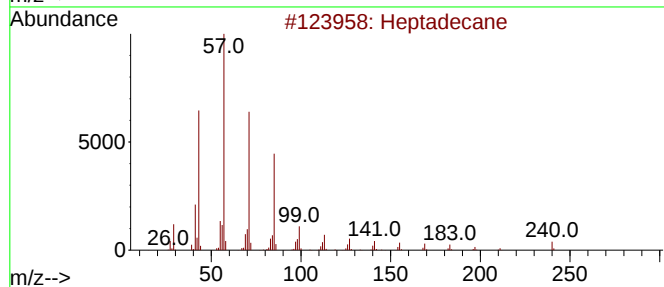
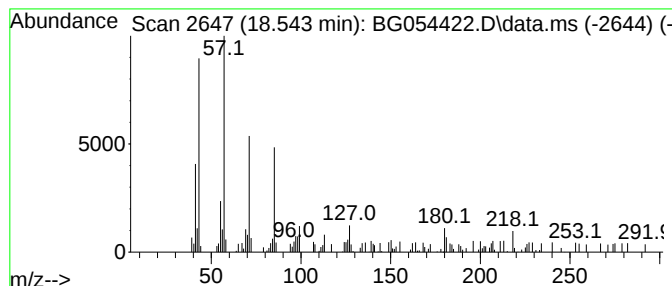
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.543	2.27 ng	34117	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	96
2		Eicosane	282	C20H42	000112-95-8	92
3		Hexadecane	226	C16H34	000544-76-3	90
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	49
5		Tetradecane	198	C14H30	000629-59-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
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Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BIN-4

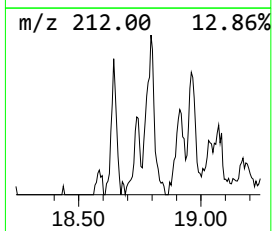
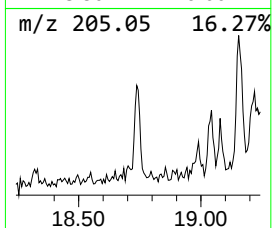
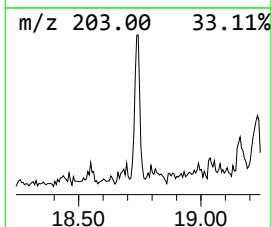
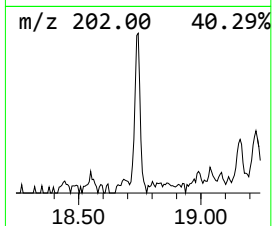
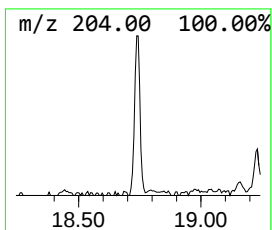
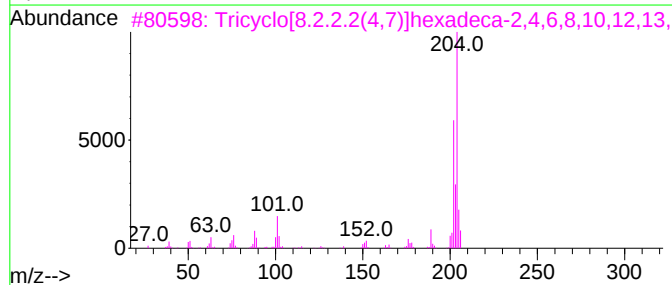
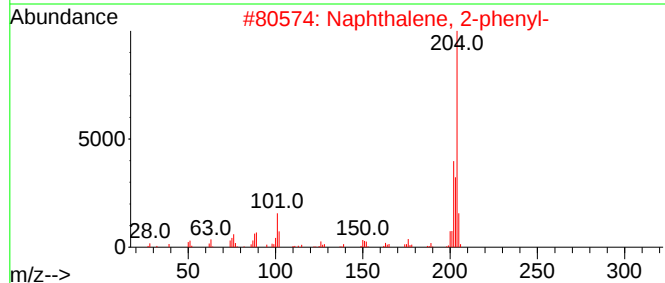
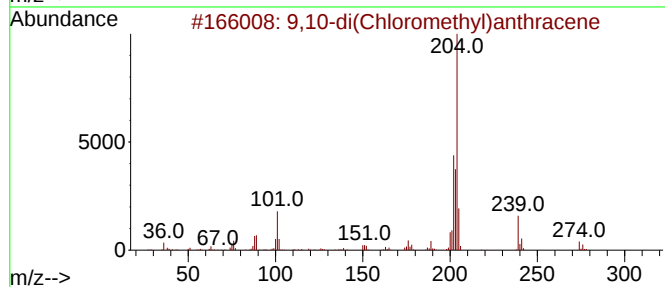
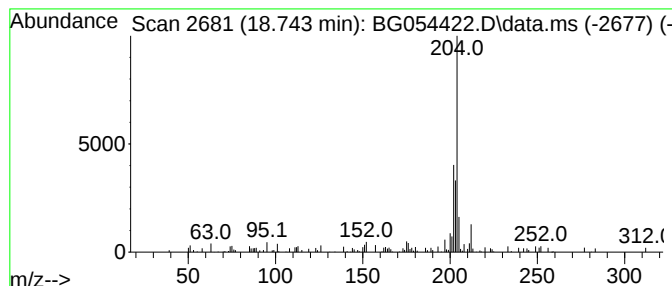
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9,10-di(Chloromethyl)anthra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.743	2.44 ng	36584	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	72
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	60
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	60
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	47
5		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
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Acq On : 27 Jul 2022 19:10
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Sample : N3892-01
Misc :
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Instrument :
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ClientSampleId :
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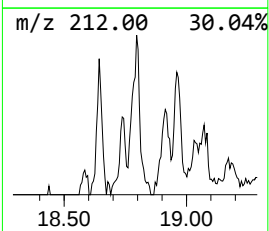
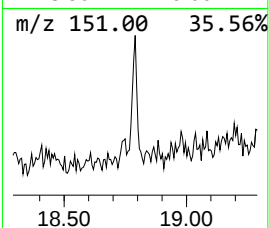
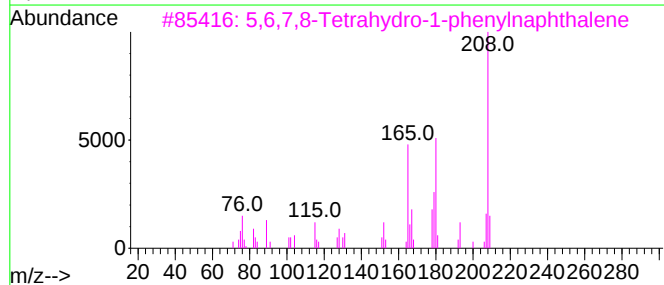
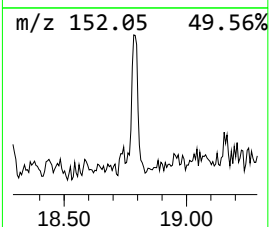
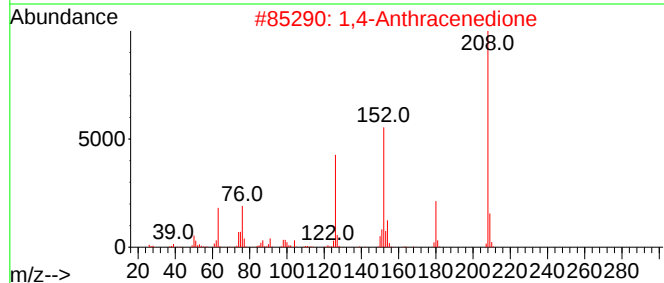
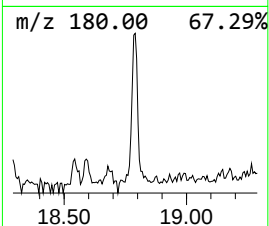
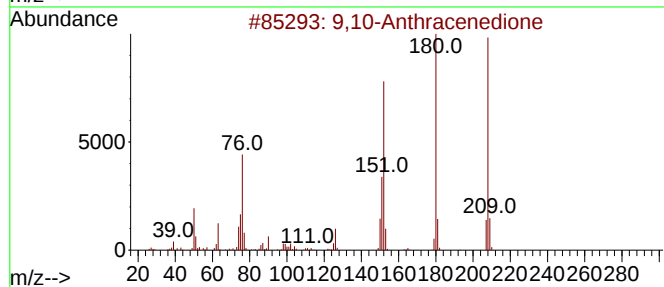
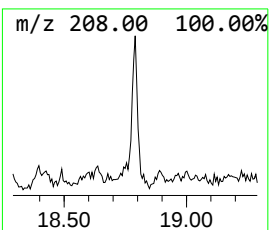
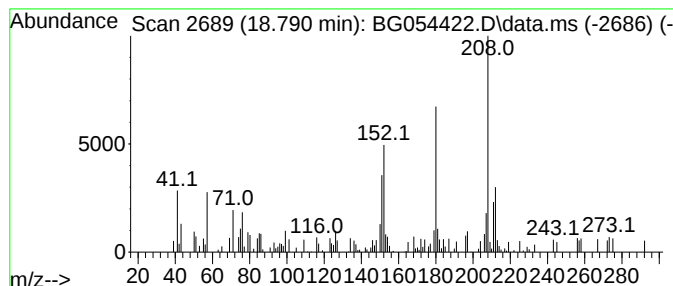
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.790	2.22 ng	33316	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	86
2		1,4-Anthracenedione	208	C14H8O2	000635-12-1	49
3		5,6,7,8-Tetrahydro-1-phenylnapht...	208	C16H16	067064-63-5	43
4		Pyrene, 1,2,3,3a,4,5-hexahydro-	208	C16H16	005385-37-5	43
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
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Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BIN-4

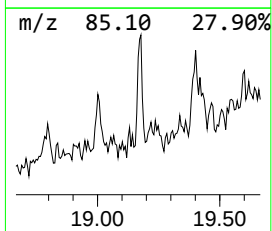
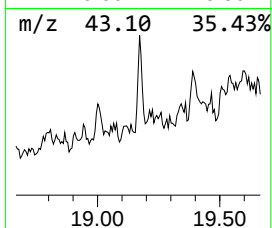
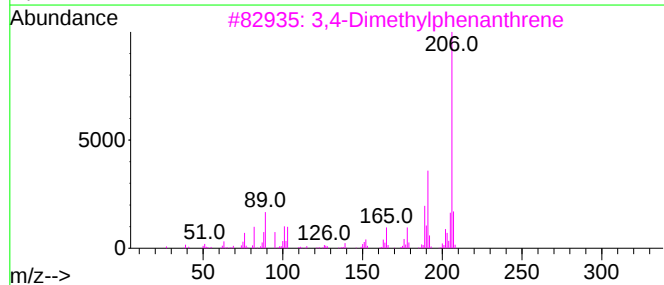
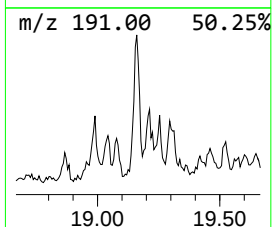
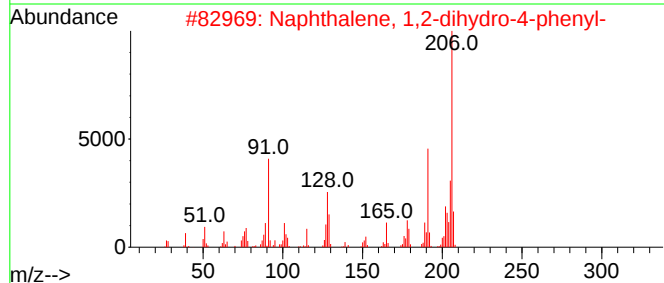
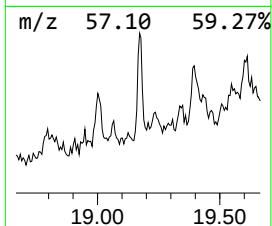
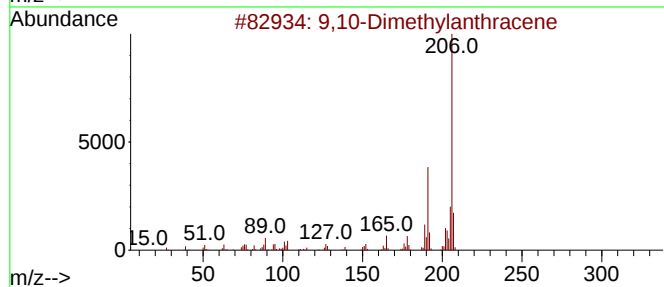
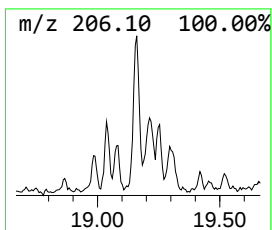
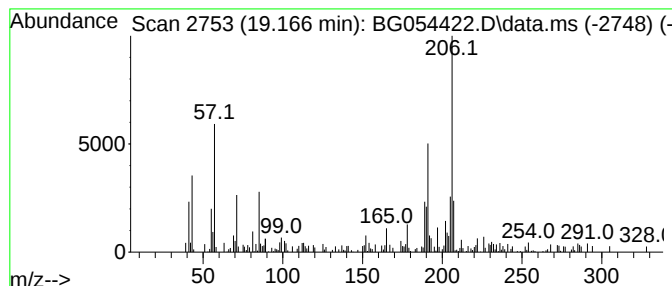
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,10-Dimethylantracene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.166	5.98 ng	89638	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	89
2			Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	64
3			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	64
4			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	62
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
Operator : CG/JU
Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BIN-4

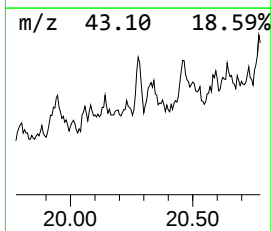
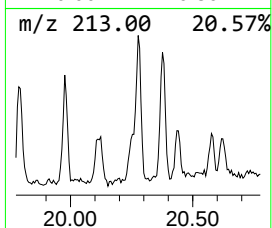
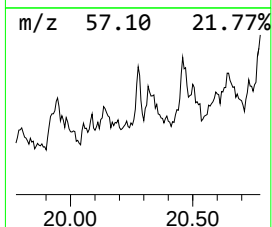
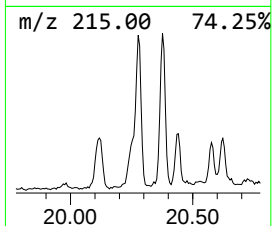
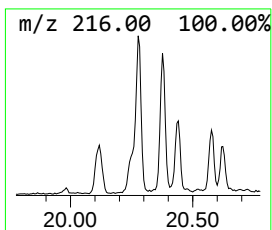
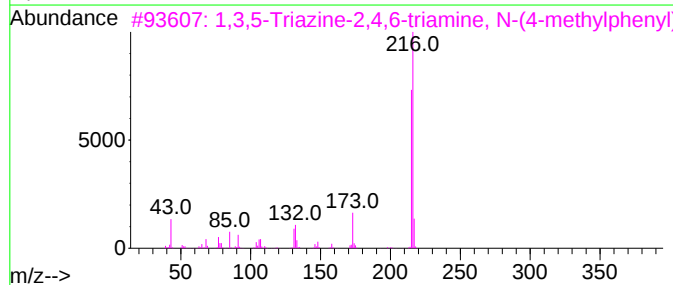
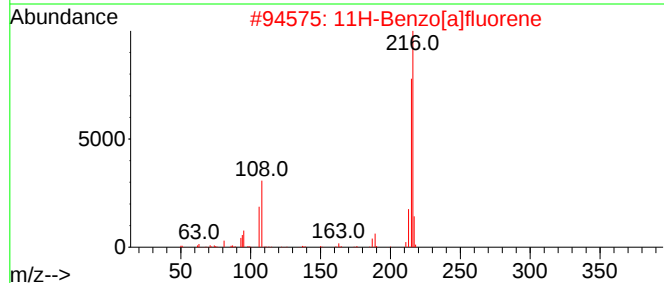
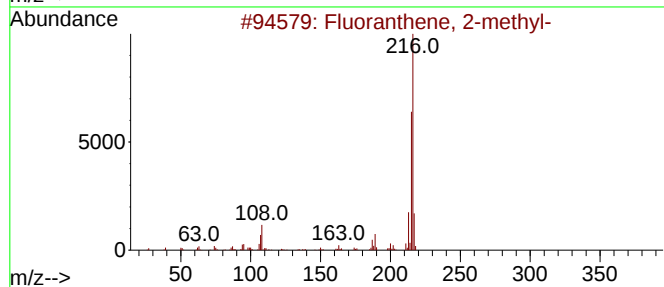
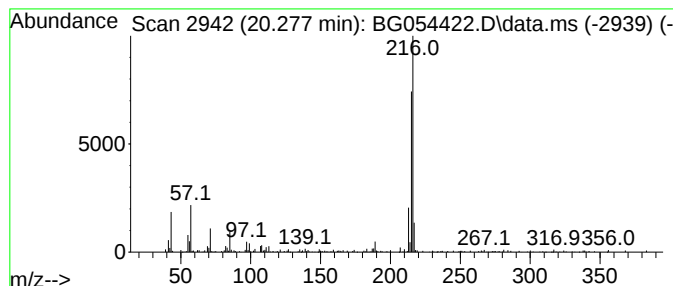
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Fluoranthene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.276	4.01 ng	150289	Chrysene-d12	21.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
3			1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	64
4			3-Trifluoromethylcinnamic acid	216	C10H7F3O2	000779-89-5	64
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072722\
Data File : BG054422.D
Acq On : 27 Jul 2022 19:10
Operator : CG/JU
Sample : N3892-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BIN-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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
Butane, 2-metho...	3.108	22.0	ng	117579	1	7.991	106960	20.0
Hexacosane	17.110	2.3	ng	34039	4	17.368	299939	20.0
Dibenzothiophene	17.186	3.5	ng	52952	4	17.368	299939	20.0
Heptadecane, 9-...	17.850	2.0	ng	30309	4	17.368	299939	20.0
Anthracene, 1-m...	18.250	6.0	ng	90408	4	17.368	299939	20.0
Anthracene, 2-m...	18.291	5.7	ng	84999	4	17.368	299939	20.0
1H-Cyclopropa[1...	18.373	2.2	ng	33140	4	17.368	299939	20.0
4H-Cyclopenta[d...	18.443	9.4	ng	141375	4	17.368	299939	20.0
Heptadecane	18.543	2.3	ng	34117	4	17.368	299939	20.0
9,10-di(Chlorom...	18.743	2.4	ng	36584	4	17.368	299939	20.0
9,10-Anthracene...	18.790	2.2	ng	33316	4	17.368	299939	20.0
9,10-Dimethylan...	19.166	6.0	ng	89638	4	17.368	299939	20.0
Fluoranthene, 2...	20.276	4.0	ng	150289	5	21.640	749441	20.0