

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
 Data File : BP009583.D
 Acq On : 29 Mar 2022 01:05
 Operator : CG/JU
 Sample : N2095-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-06-4.0-5.0

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_P\Methods\8270E-BP031722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009583.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	15	rVB	447329	590668	3.24%	0.565%
2	5.375	399	406	429	rBV	4552931	7977981	43.73%	7.630%
3	6.958	664	675	705	rBV	4260332	8210529	45.01%	7.853%
4	7.769	805	813	821	rBV	919402	1662428	9.11%	1.590%
5	8.922	995	1009	1029	rBV	2836043	5384335	29.52%	5.150%
6	10.557	1276	1287	1303	rBV	1175227	2352296	12.90%	2.250%
7	13.034	1700	1708	1725	rBV	7778096	12652845	69.36%	12.101%
8	14.410	1935	1942	1949	rBV	1872794	3155667	17.30%	3.018%
9	14.475	1949	1953	1963	rVB	129418	241273	1.32%	0.231%
10	15.916	2191	2198	2222	rBV	6009304	10101657	55.38%	9.661%
11	16.151	2234	2238	2247	rVB	143716	264589	1.45%	0.253%
12	16.951	2367	2374	2378	rVV3	115144	251609	1.38%	0.241%
13	16.998	2378	2382	2393	rVB2	183603	354549	1.94%	0.339%
14	17.175	2406	2412	2416	rBV	2404637	3762250	20.62%	3.598%
15	17.222	2416	2420	2426	rVB	1902189	2847159	15.61%	2.723%
16	17.310	2431	2435	2441	rVB	289683	481889	2.64%	0.461%
17	17.698	2497	2501	2507	rVB2	201059	272535	1.49%	0.261%
18	17.763	2508	2512	2516	rBV2	166022	251092	1.38%	0.240%
19	18.051	2556	2561	2564	rBV	591416	878043	4.81%	0.840%
20	18.092	2564	2568	2576	rVB2	1014357	1842242	10.10%	1.762%
21	18.239	2588	2593	2597	rBV2	649601	1209605	6.63%	1.157%
22	18.275	2597	2599	2605	rVB	426288	554517	3.04%	0.530%
23	18.369	2611	2615	2622	rBV2	245733	413976	2.27%	0.396%
24	18.539	2640	2644	2647	rBV	374045	506026	2.77%	0.484%
25	18.586	2648	2652	2660	rVB3	277488	528163	2.90%	0.505%
26	18.751	2677	2680	2682	rBV2	155256	201031	1.10%	0.192%
27	18.828	2690	2693	2697	rBV2	156792	223330	1.22%	0.214%
28	18.945	2709	2713	2717	rBV	408963	658966	3.61%	0.630%
29	18.986	2717	2720	2726	rVB4	388628	642740	3.52%	0.615%
30	19.086	2733	2737	2743	rBV4	261617	513631	2.82%	0.491%
31	19.204	2752	2757	2764	rBV	2279000	3210946	17.60%	3.071%
32	19.557	2810	2817	2821	rBV	3130047	5107017	28.00%	4.884%
33	19.757	2846	2851	2862	rVB	13371735	18241700	100.00%	17.447%
34	20.557	2984	2987	2991	rBV	591189	696153	3.82%	0.666%
35	21.022	3063	3066	3073	rVB4	1055402	1964235	10.77%	1.879%
36	21.292	3106	3112	3116	rBV2	3591490	6349859	34.81%	6.073%

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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

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_P\Methods\8270E-BP031722.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

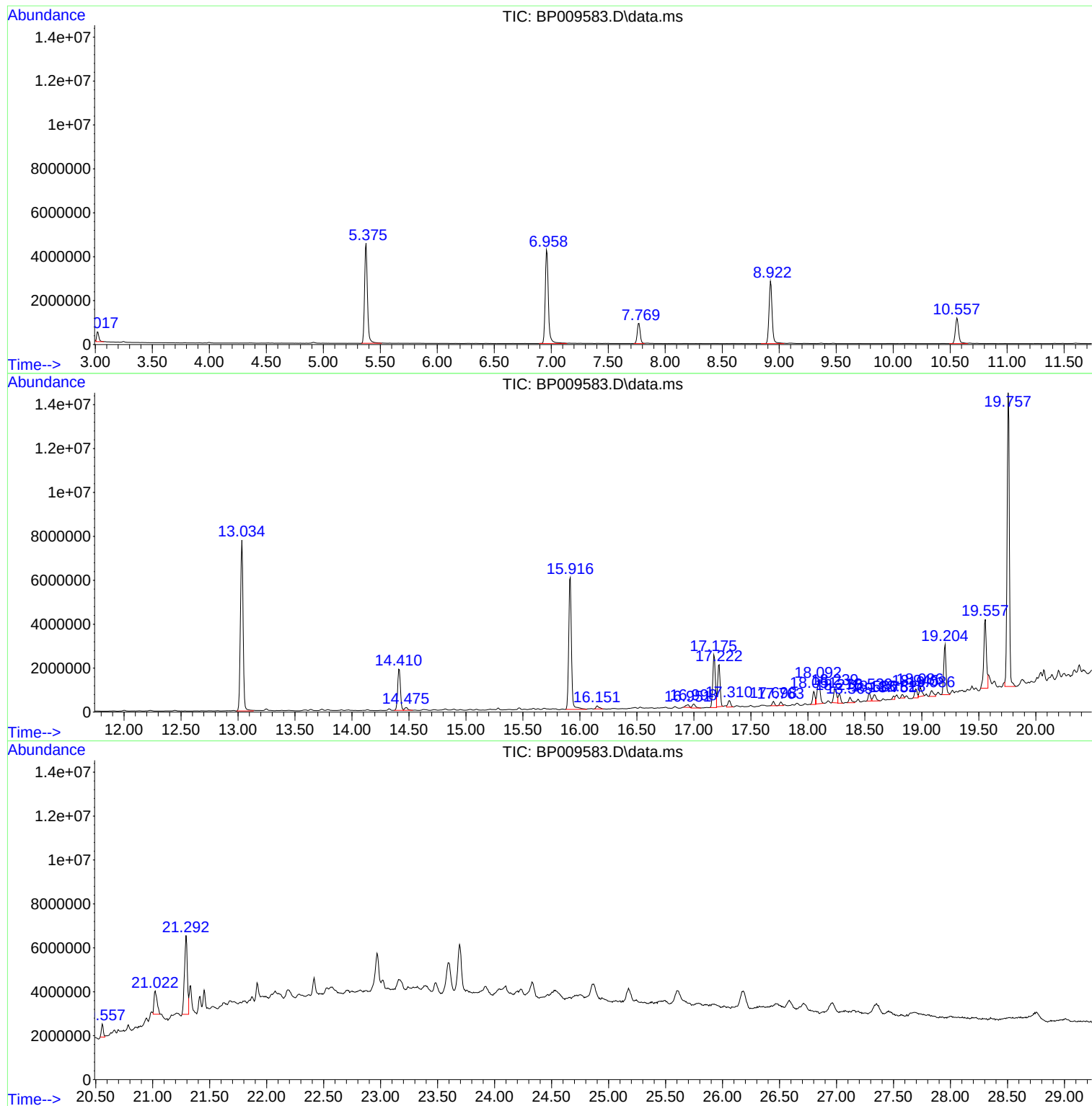
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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ClientSampleId :
SB-06-4.0-5.0

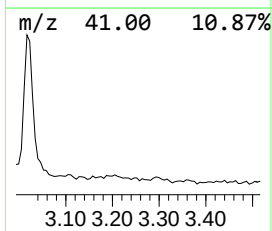
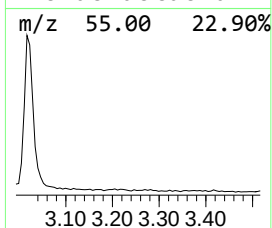
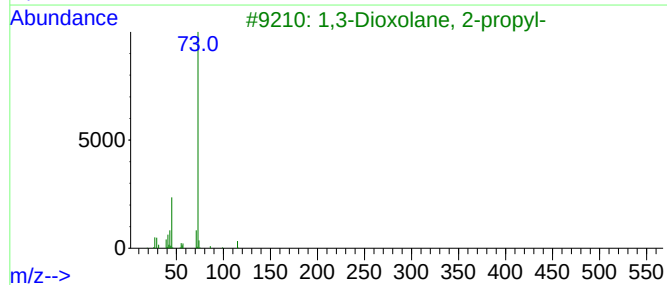
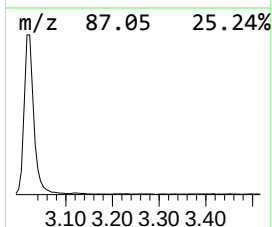
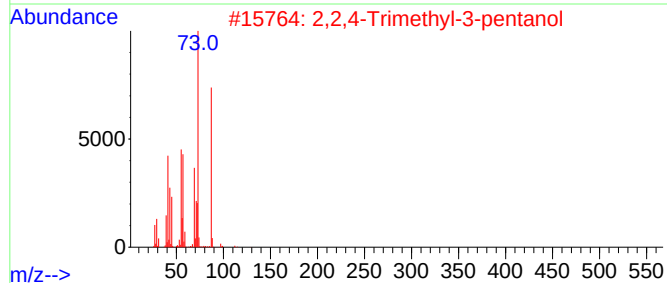
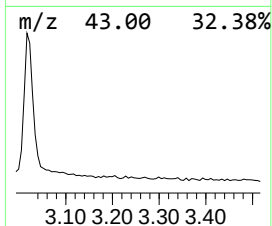
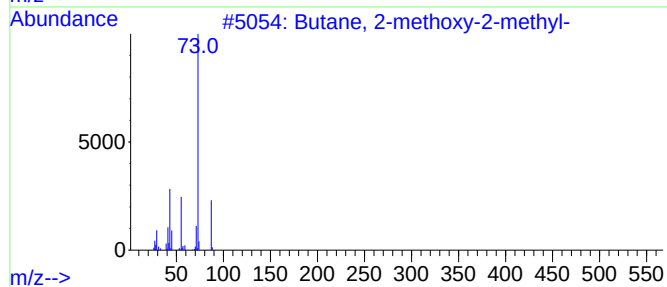
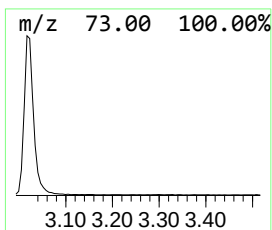
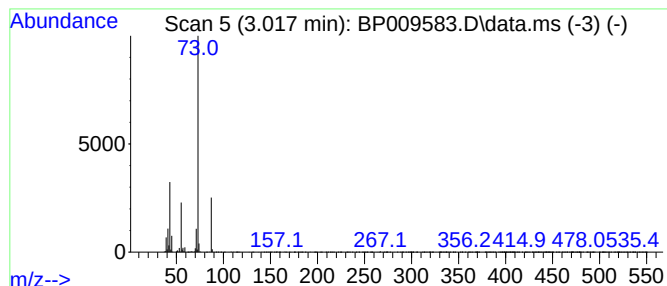
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	7.11 ng	590668	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

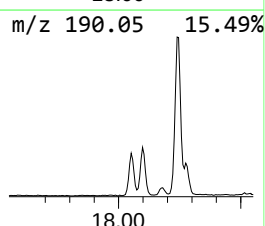
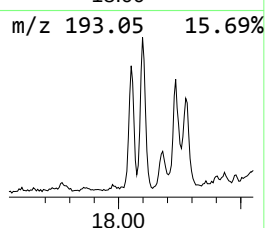
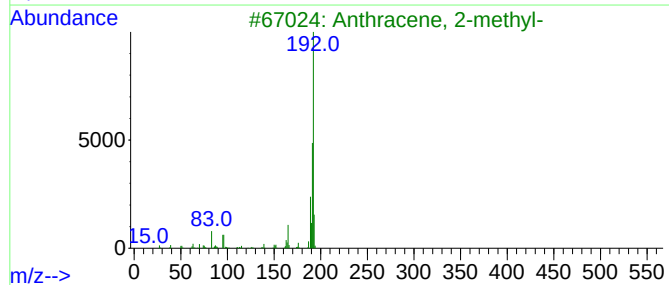
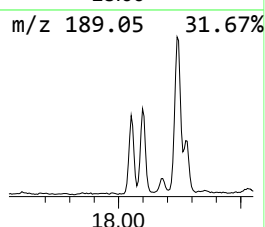
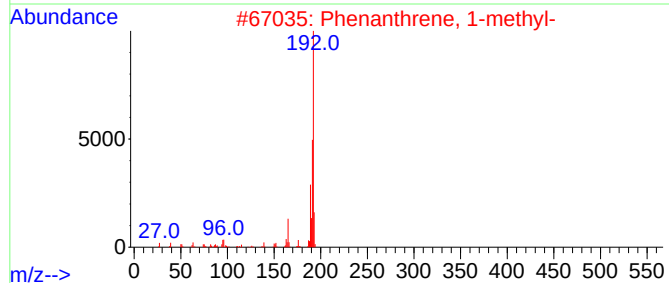
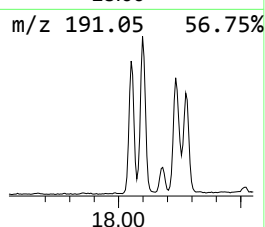
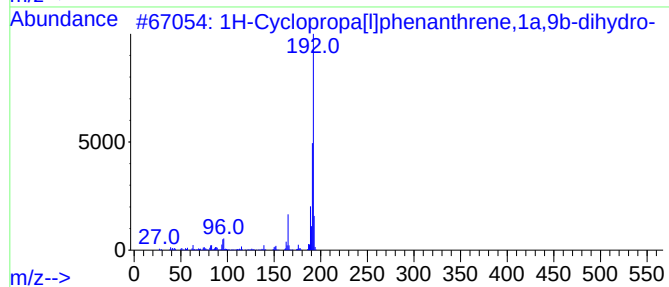
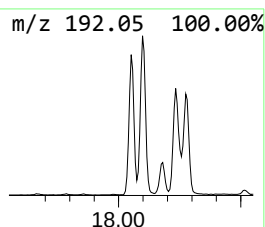
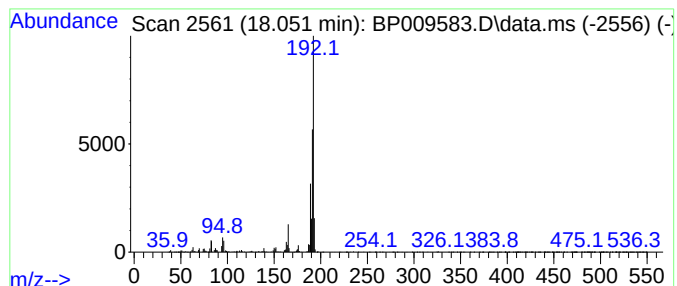
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	4.67 ng	878043	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95



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Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

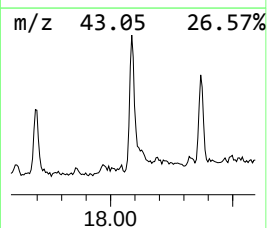
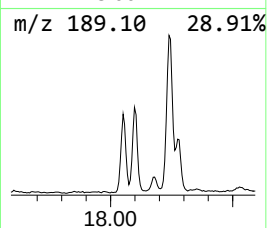
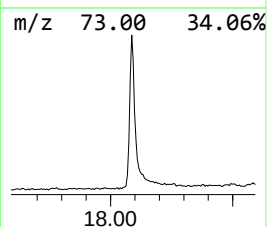
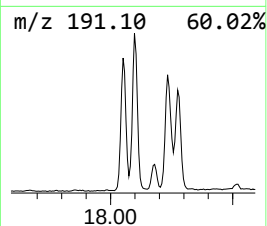
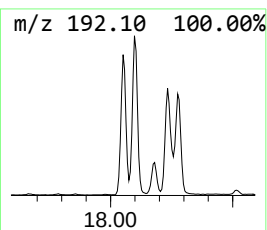
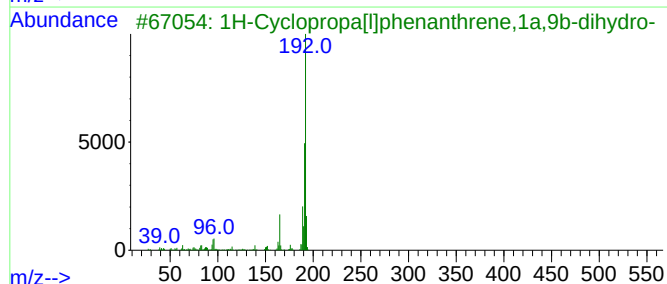
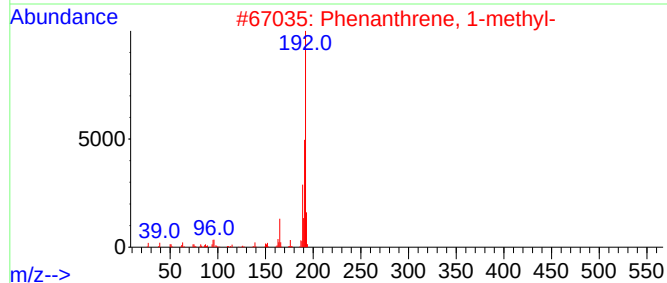
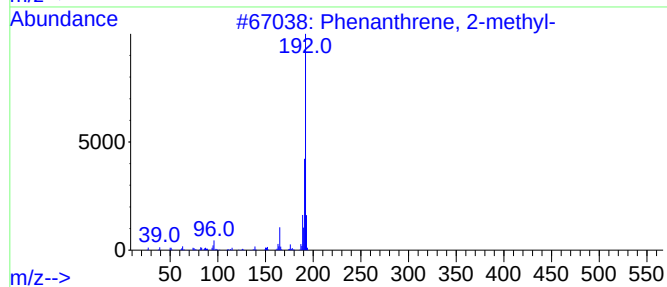
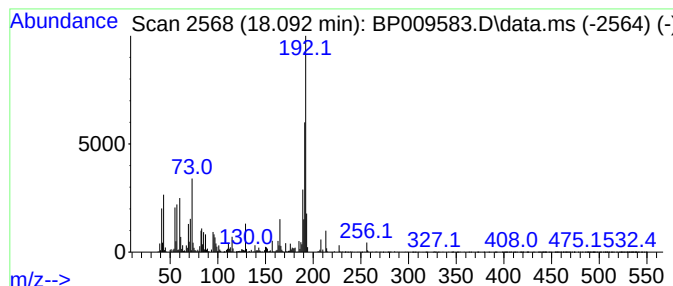
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	9.79 ng	1842240	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	92



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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

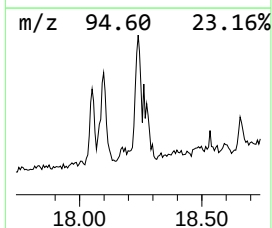
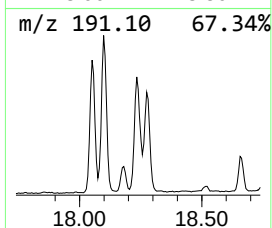
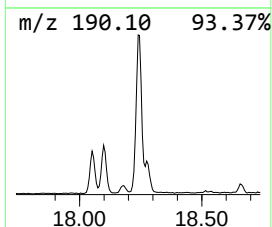
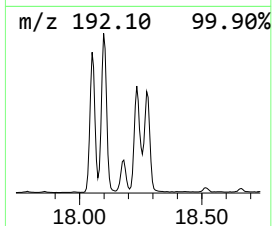
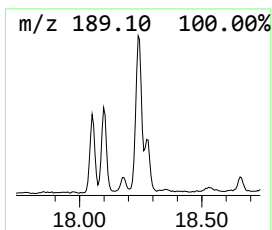
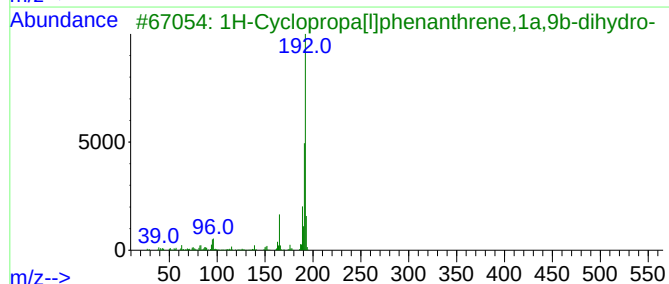
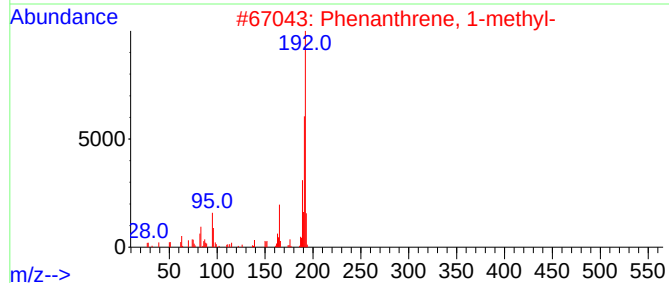
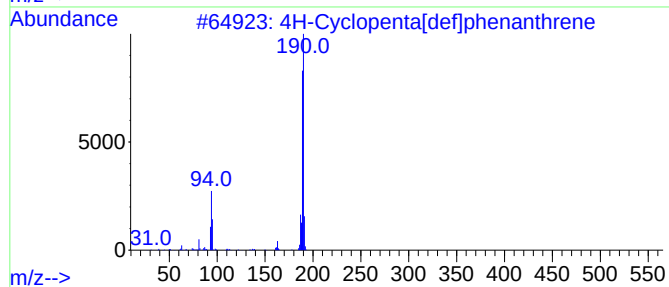
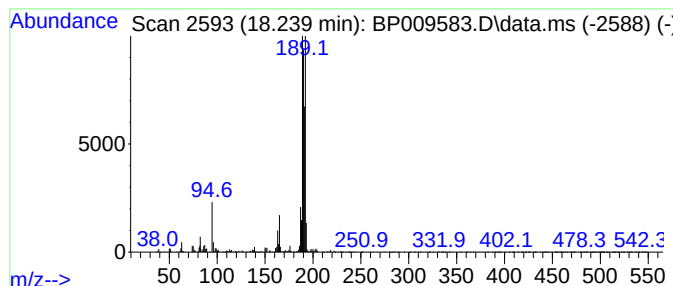
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	6.43 ng	1209610	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	42
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	42



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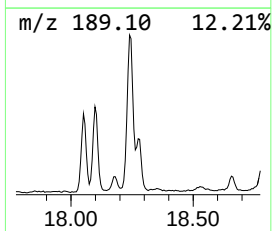
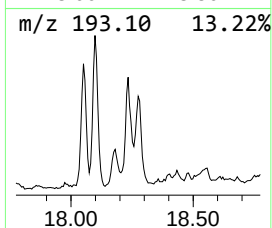
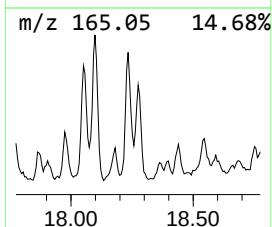
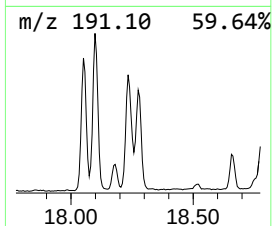
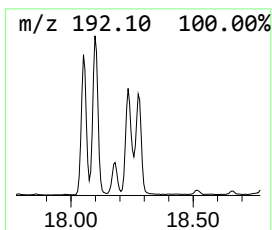
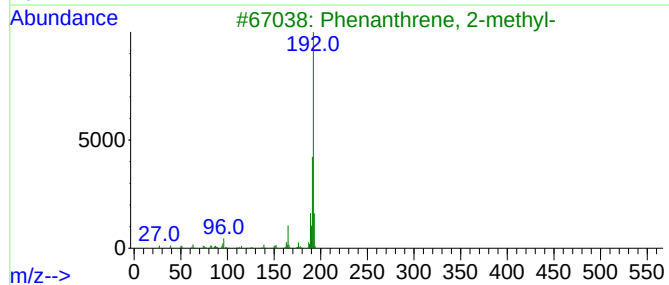
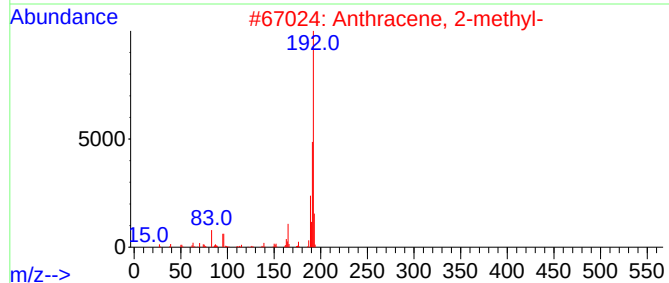
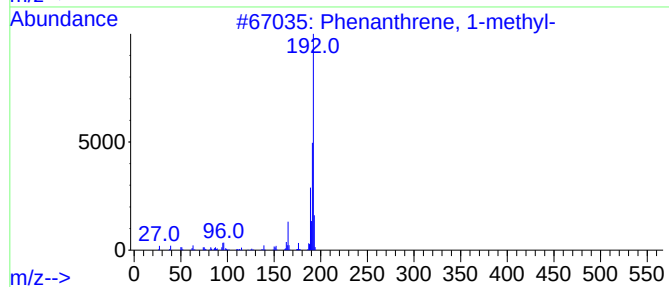
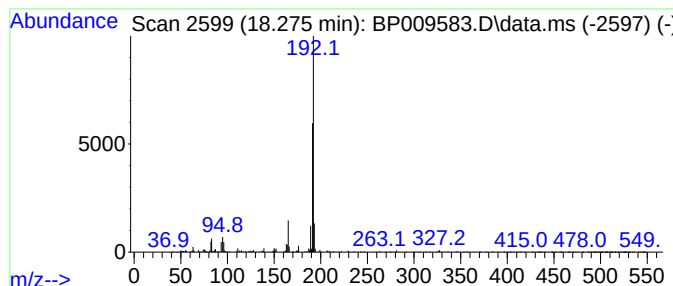
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.275	2.95 ng	554517	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
5			1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	90



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Operator : CG/JU
Sample : N2095-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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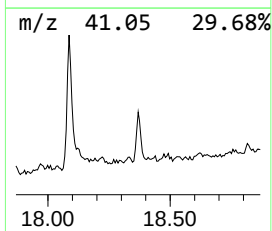
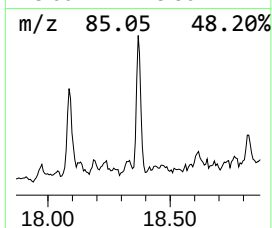
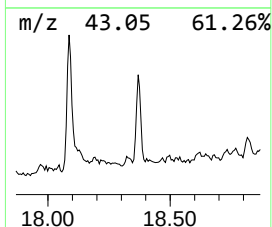
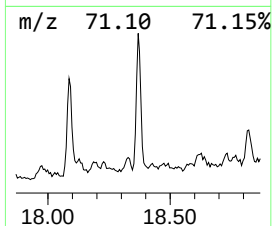
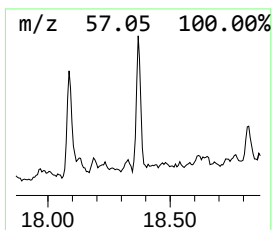
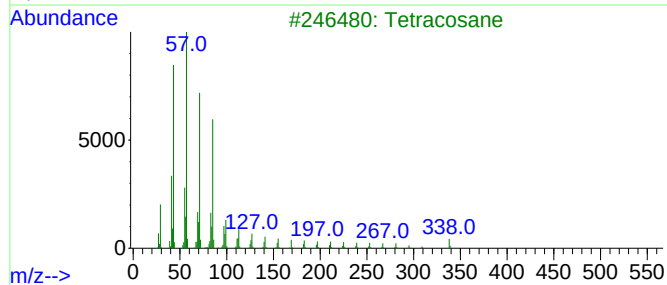
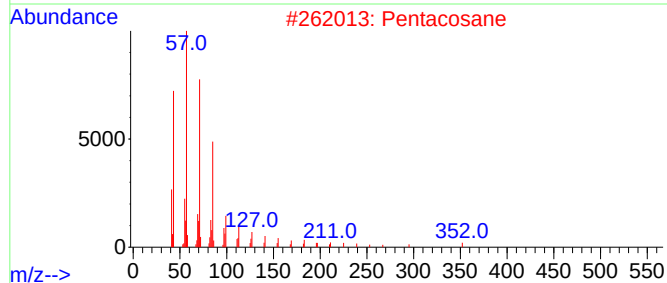
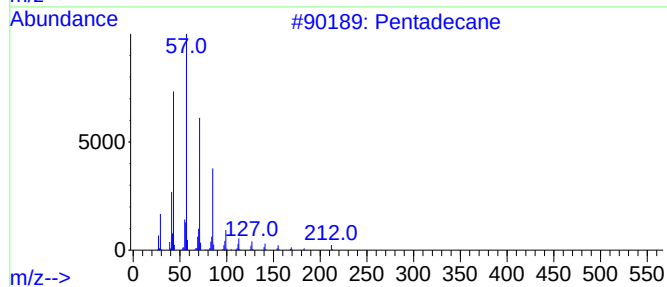
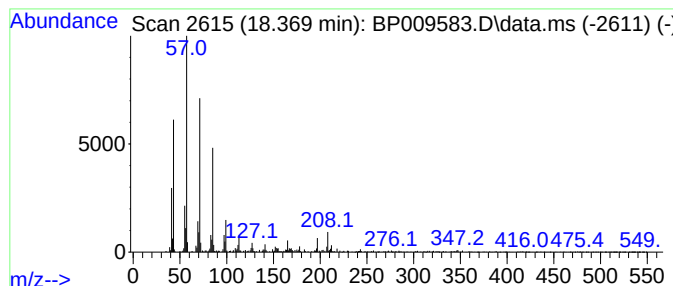
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	2.20 ng	413976	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	93
2			Pentacosane	352	C25H52	000629-99-2	90
3			Tetracosane	338	C24H50	000646-31-1	81
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	74
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
Operator : CG/JU
Sample : N2095-06
Misc :
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ClientSampleId :
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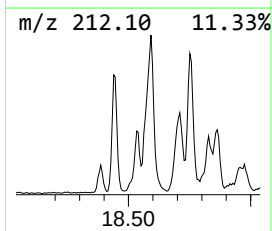
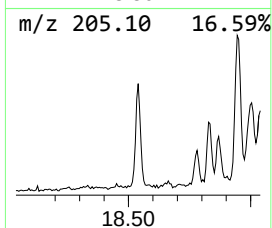
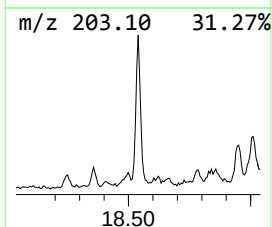
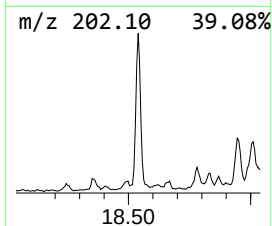
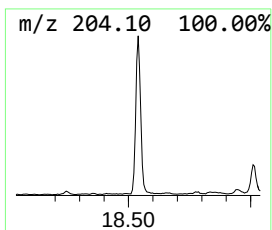
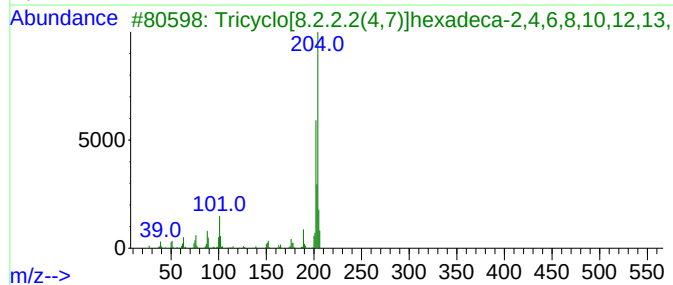
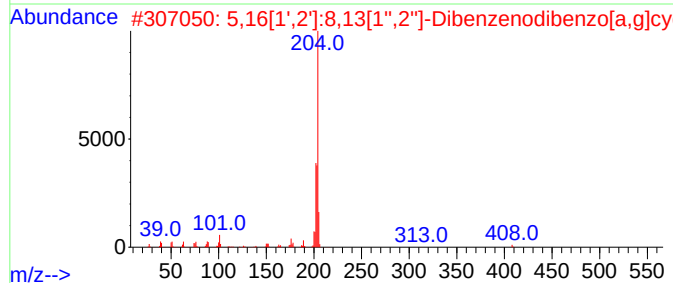
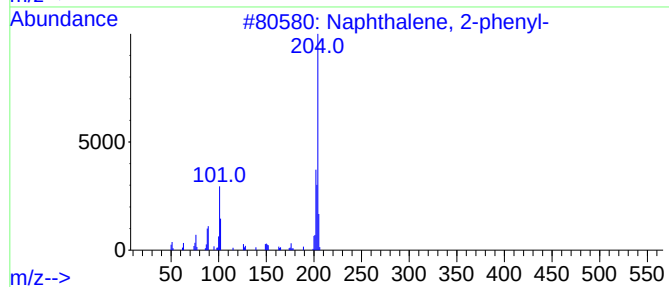
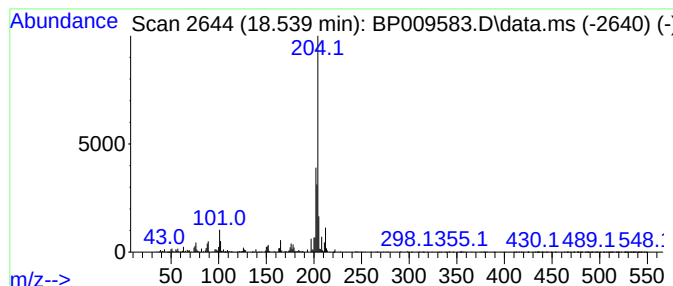
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	2.69 ng	506026	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	83
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
4			9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58
5			5H-Dibenzo[a,d]cycloheptene, 5-m...	204	C16H12	002975-79-3	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
Operator : CG/JU
Sample : N2095-06
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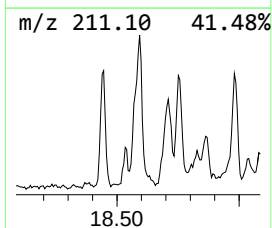
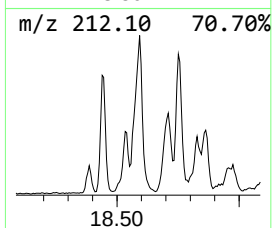
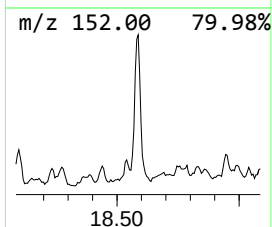
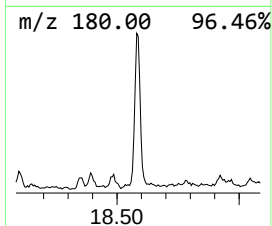
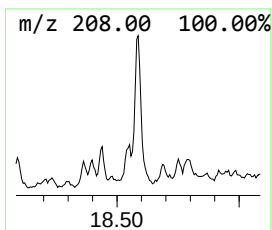
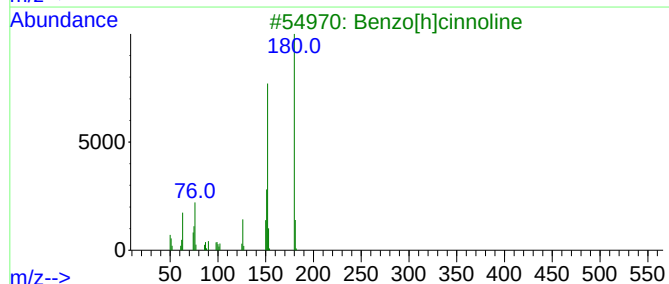
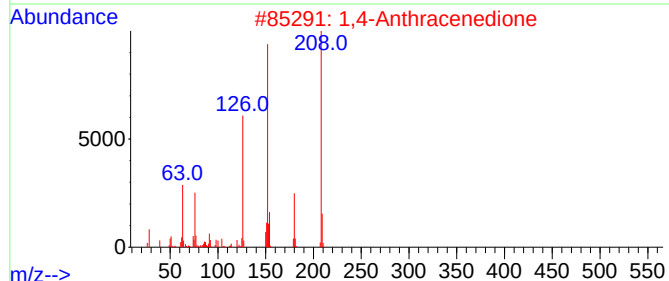
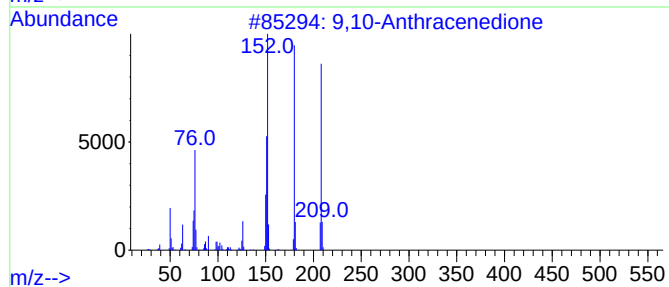
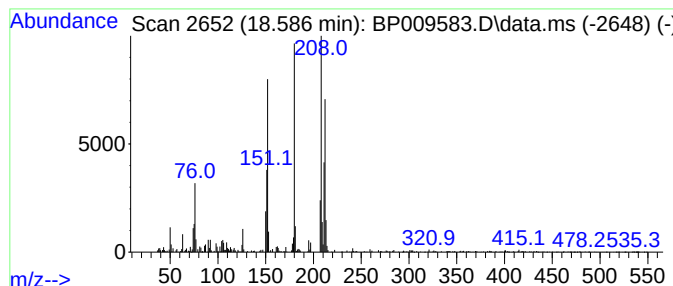
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SB-06-4.0-5.0

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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 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.586	2.81 ng	528163	Phenanthrene-d10		17.175	
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione		208	C14H8O2	000084-65-1	97
2	1,4-Anthracenedione		208	C14H8O2	000635-12-1	50
3	Benzo[h]cinnoline		180	C12H8N2	000230-31-9	50
4	Benzo[c]cinnoline		180	C12H8N2	000230-17-1	50
5	1H-Phenalen-1-one		180	C13H8O	000548-39-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
Operator : CG/JU
Sample : N2095-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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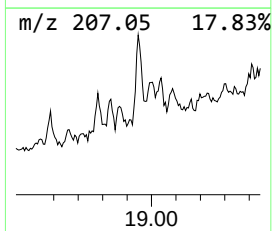
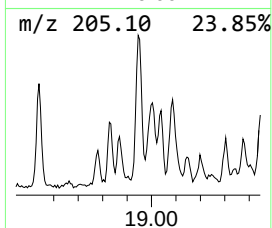
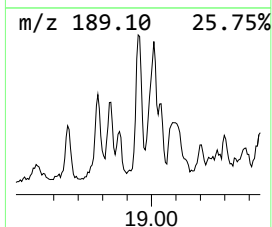
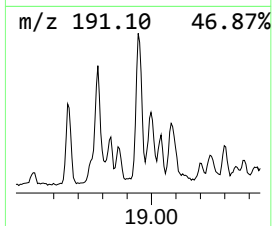
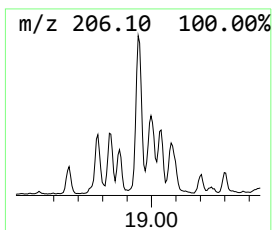
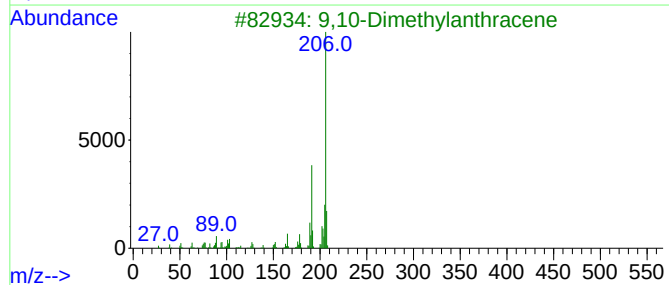
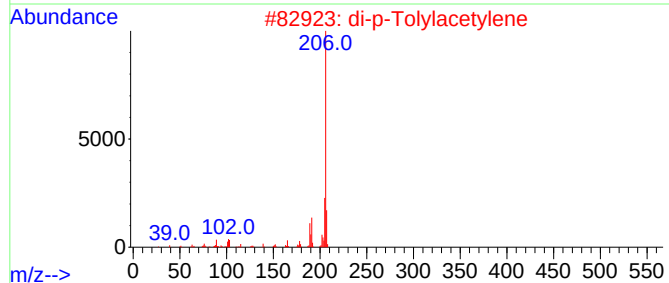
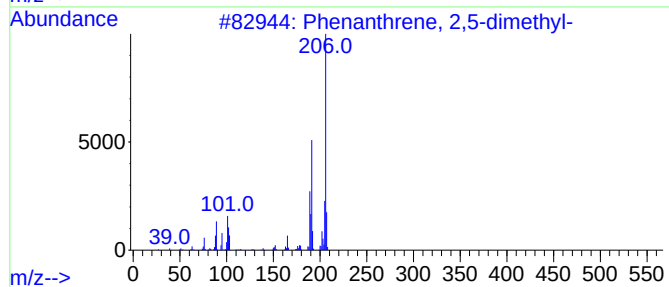
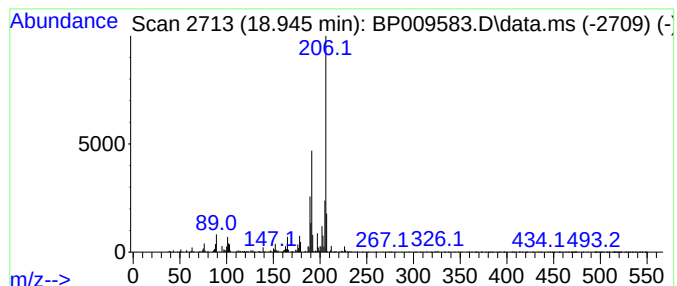
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.945	3.50 ng	658966	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	91
3			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
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ALS Vial : 17 Sample Multiplier: 1

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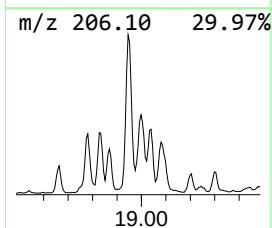
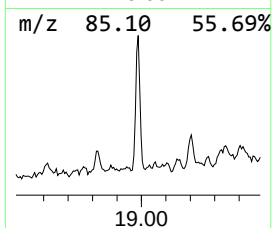
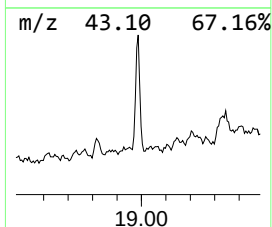
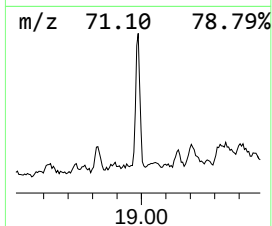
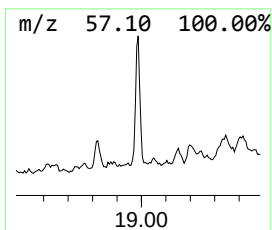
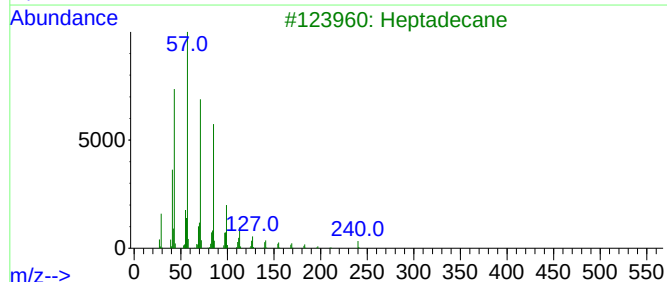
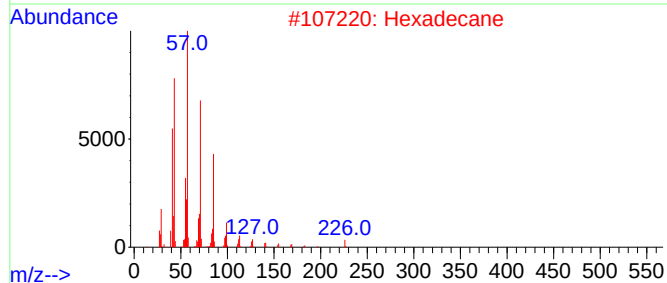
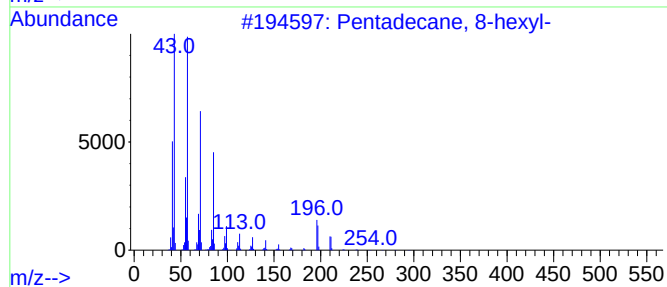
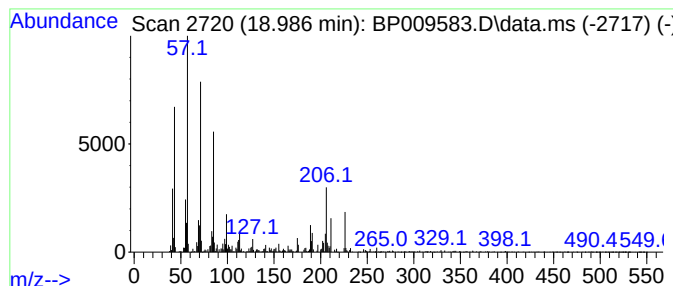
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane, 8-hexyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.986	3.42 ng	642740	Phenanthrene-d10	17.175

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	76
2		Hexadecane	226	C16H34	000544-76-3	74
3		Heptadecane	240	C17H36	000629-78-7	68
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	68
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
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Acq On : 29 Mar 2022 01:05
Operator : CG/JU
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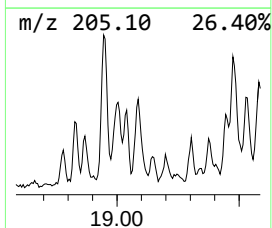
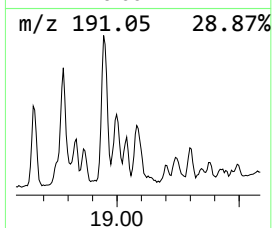
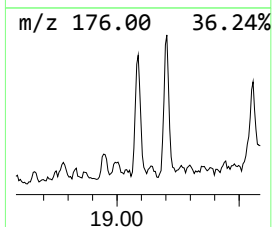
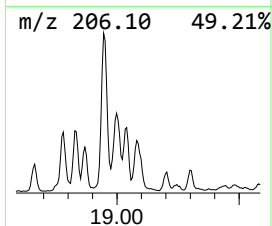
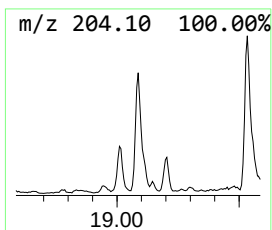
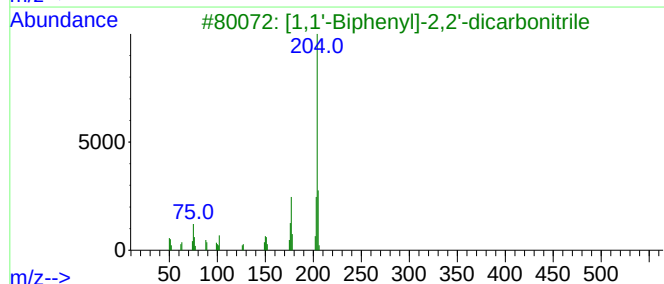
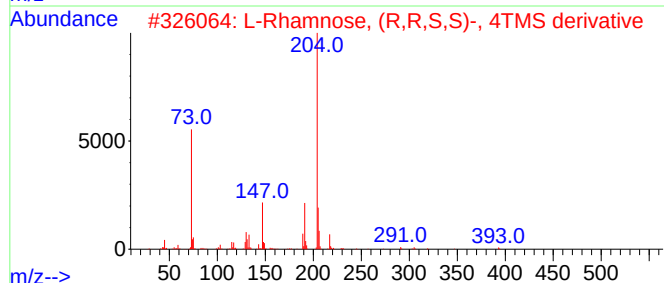
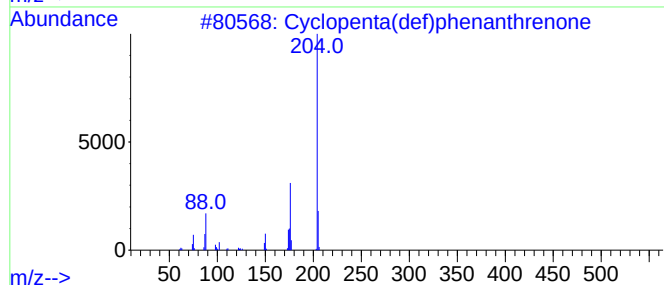
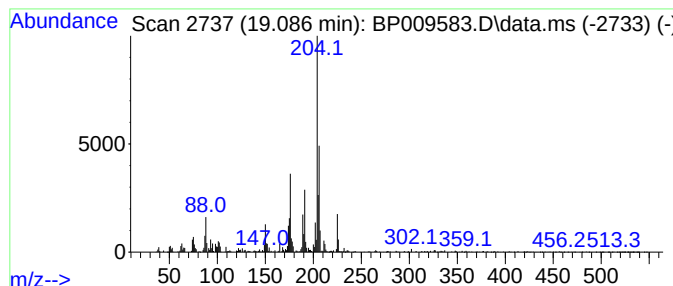
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	2.73 ng	513631	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	89
2			L-Rhamnose, (R,R,S,S)-, 4TMS der...	452	C18H44O5Si4	108392-01-4	47
3			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	43
4			[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	40
5			(S)-N-(Furan-2-ylmethyl)-1-(1,2,...	511	C27H41N3O3Si2	1010498-97-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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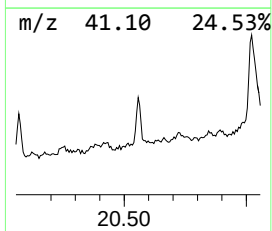
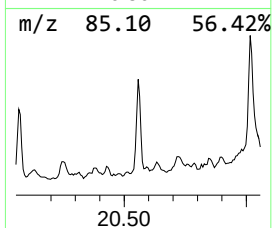
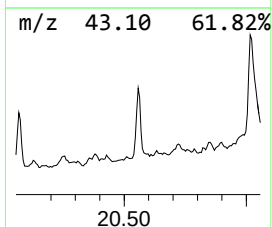
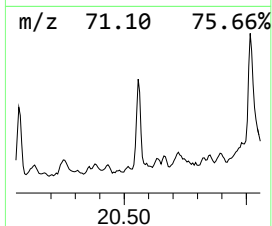
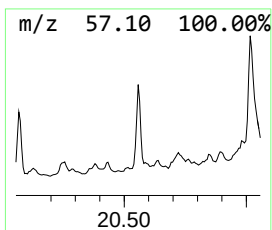
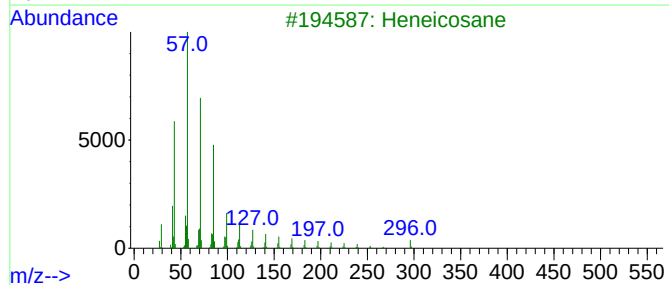
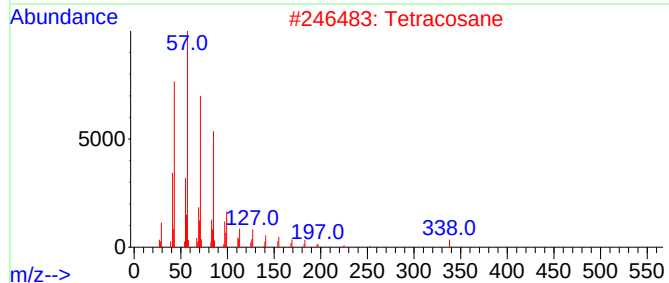
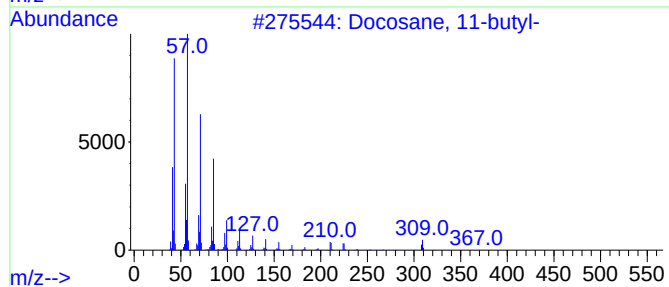
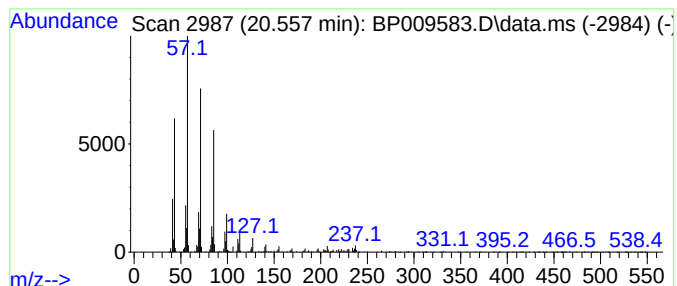
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane, 11-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.557	2.19 ng	696153	Chrysene-d12	21.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
2		Tetracosane	338	C24H50	000646-31-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
5		Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
Operator : CG/JU
Sample : N2095-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

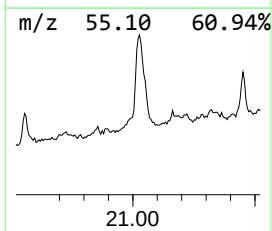
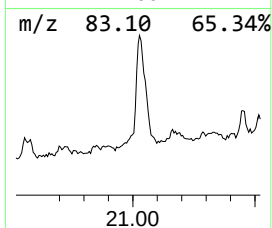
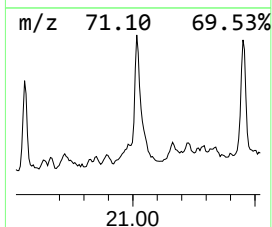
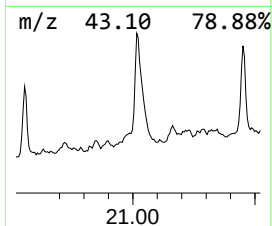
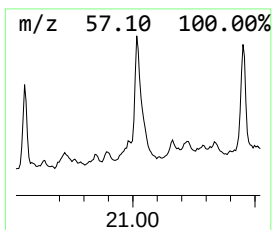
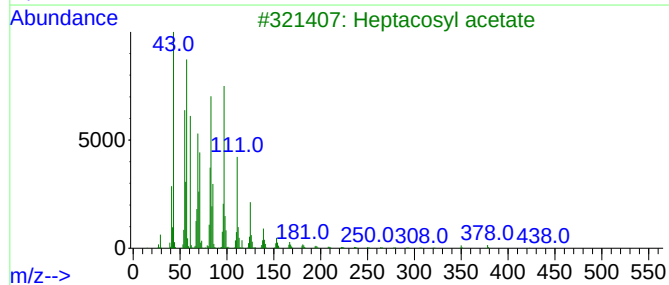
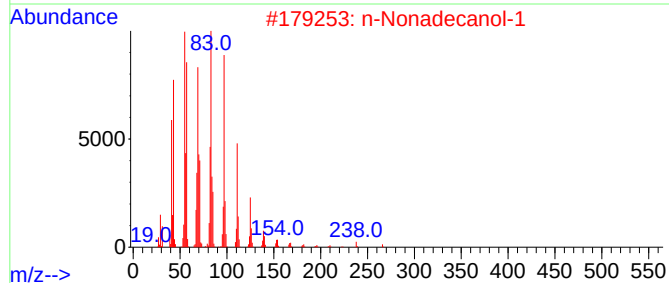
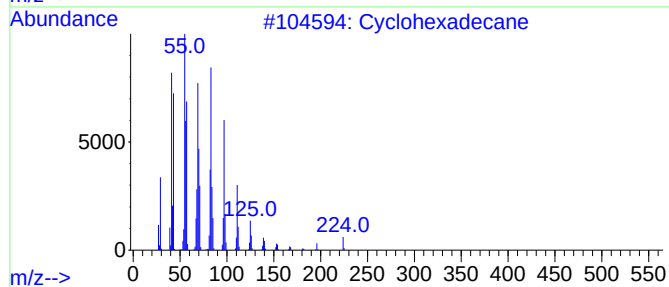
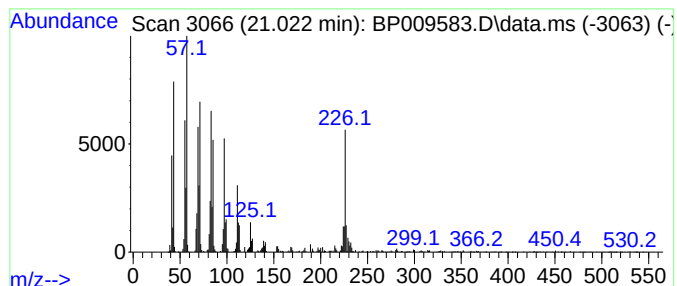
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclohexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	6.19 ng	1964240	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	97
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	87
3			Heptacosyl acetate	438	C29H58O2	1000351-78-2	86
4			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	83
5			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032822\
Data File : BP009583.D
Acq On : 29 Mar 2022 01:05
Operator : CG/JU
Sample : N2095-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-06-4.0-5.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.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...	3.017	7.1	ng	590668	1	7.769	1662430	20.0
1H-Cyclopropa[1...	18.051	4.7	ng	878043	4	17.175	3762250	20.0
Phenanthrene, 2...	18.092	9.8	ng	1842240	4	17.175	3762250	20.0
4H-Cyclopenta[d...	18.239	6.4	ng	1209610	4	17.175	3762250	20.0
Phenanthrene, 1...	18.275	3.0	ng	554517	4	17.175	3762250	20.0
Pentadecane	18.369	2.2	ng	413976	4	17.175	3762250	20.0
Naphthalene, 2-...	18.539	2.7	ng	506026	4	17.175	3762250	20.0
9,10-Anthracene...	18.586	2.8	ng	528163	4	17.175	3762250	20.0
Phenanthrene, 2...	18.945	3.5	ng	658966	4	17.175	3762250	20.0
Pentadecane, 8-...	18.986	3.4	ng	642740	4	17.175	3762250	20.0
Cyclopenta(def)...	19.086	2.7	ng	513631	4	17.175	3762250	20.0
Docosane, 11-bu...	20.557	2.2	ng	696153	5	21.292	6349860	20.0
Cyclohexadecane	21.022	6.2	ng	1964240	5	21.292	6349860	20.0