

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
 Data File : BP003228.D
 Acq On : 08 Sep 2020 15:48
 Operator : CG/JU
 Sample : L3916-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP082420.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	32	rVB	1136189	1241960	20.77%	2.024%
2	4.793	318	324	331	rBV	48644	69103	1.16%	0.113%
3	5.264	397	404	421	rBV	2008189	2887533	48.29%	4.707%
4	6.840	661	672	685	rBV	1647638	2716009	45.42%	4.427%
5	7.652	803	810	818	rBV	709529	1111993	18.60%	1.813%
6	7.734	818	824	847	rVV	141589	350960	5.87%	0.572%
7	8.799	992	1005	1019	rBV	1236260	2091965	34.98%	3.410%
8	10.428	1275	1282	1289	rBV	841588	1430945	23.93%	2.332%
9	11.887	1524	1530	1536	rBV	63321	122574	2.05%	0.200%
10	12.910	1697	1704	1724	rBV	2942294	4580523	76.60%	7.466%
11	13.622	1819	1825	1830	rBV4	84130	148785	2.49%	0.243%
12	13.751	1841	1847	1856	rBV	204387	327364	5.47%	0.534%
13	14.016	1886	1892	1899	rBV	323384	503478	8.42%	0.821%
14	14.151	1908	1915	1925	rBV2	100080	279438	4.67%	0.455%
15	14.293	1933	1939	1946	rBV	1106902	1650126	27.59%	2.690%
16	14.357	1946	1950	1959	rVB	135984	221051	3.70%	0.360%
17	14.693	2003	2007	2012	rBV2	65457	107905	1.80%	0.176%
18	15.345	2113	2118	2130	rVB	167439	278934	4.66%	0.455%
19	15.657	2165	2171	2180	rVB4	67347	190528	3.19%	0.311%
20	15.787	2188	2193	2203	rBV	1976264	3026551	50.61%	4.933%
21	16.145	2250	2254	2258	rBV	68767	102073	1.71%	0.166%
22	16.204	2261	2264	2269	rVV2	63503	109581	1.83%	0.179%
23	16.263	2269	2274	2279	rVV3	107145	175629	2.94%	0.286%
24	16.357	2286	2290	2294	rBV3	45321	69128	1.16%	0.113%
25	16.469	2303	2309	2313	rBV4	78154	169067	2.83%	0.276%
26	16.704	2345	2349	2357	rVB3	61349	104205	1.74%	0.170%
27	16.863	2373	2376	2388	rVB	118189	225573	3.77%	0.368%
28	16.969	2391	2394	2400	rBV6	51844	84340	1.41%	0.137%
29	17.045	2401	2407	2411	rBV	1361493	2062106	34.48%	3.361%
30	17.087	2411	2414	2423	rVV	2048934	2890767	48.34%	4.712%
31	17.175	2424	2429	2435	rVB	542219	846052	14.15%	1.379%
32	17.245	2437	2441	2447	rVV2	97528	159599	2.67%	0.260%
33	17.451	2471	2476	2481	rBV	186478	276450	4.62%	0.451%
34	17.634	2504	2507	2511	rVB2	64089	78226	1.31%	0.128%

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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

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.922	2552	2556	2559	rBV	284628	355924	5.95%	0.580%
36	17.969	2559	2564	2572	rBV	1591909	2257616	37.75%	3.680%
37	18.045	2574	2577	2582	rVB	160914	193743	3.24%	0.316%
38	18.110	2583	2588	2592	rBV2	536492	902419	15.09%	1.471%
39	18.410	2635	2639	2642	rBV	259904	362404	6.06%	0.591%
40	18.445	2642	2645	2652	rVB2	272309	433929	7.26%	0.707%
41	18.698	2686	2688	2691	rBV3	187485	227404	3.80%	0.371%
42	19.069	2747	2751	2757	rVB	4247013	5503060	92.02%	8.970%
43	19.210	2772	2775	2779	rBV2	875716	1129400	18.89%	1.841%
44	19.422	2803	2811	2818	rBV	3519447	5980059	100.00%	9.747%
45	19.622	2838	2845	2849	rBV	4007931	5947116	99.45%	9.694%
46	19.904	2891	2893	2898	rVB	787514	909443	15.21%	1.482%
47	21.133	3099	3102	3107	rBV3	2094559	4068573	68.04%	6.632%
48	21.180	3108	3110	3117	rVB	2051422	2388468	39.94%	3.893%

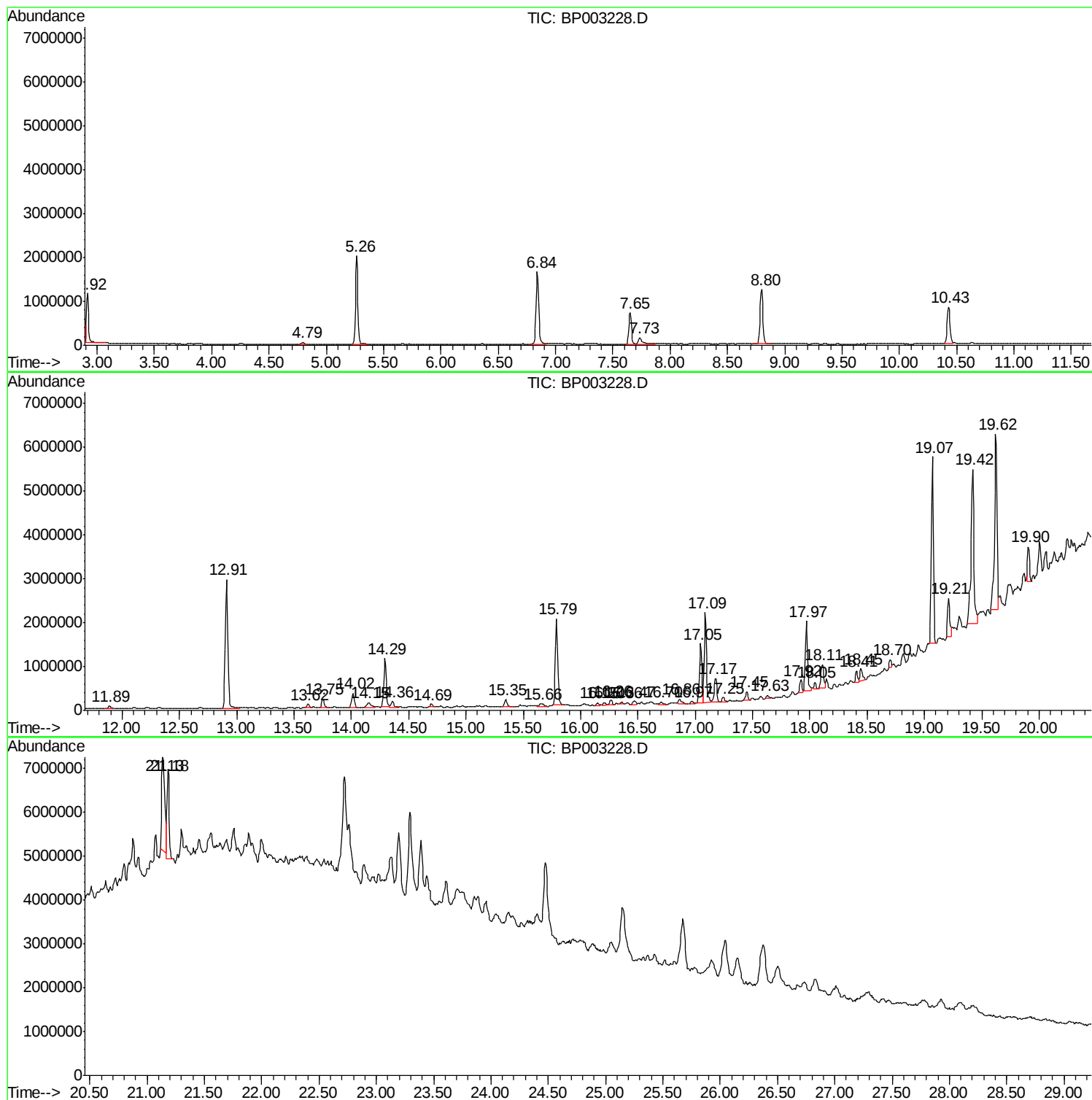
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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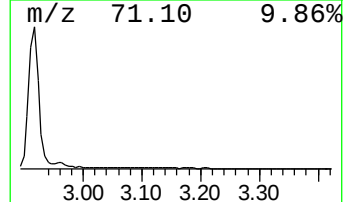
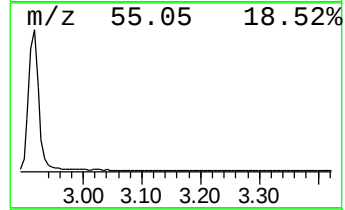
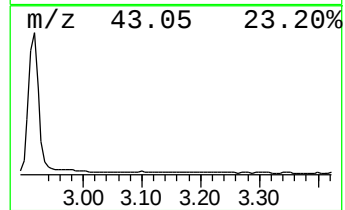
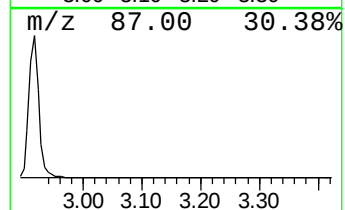
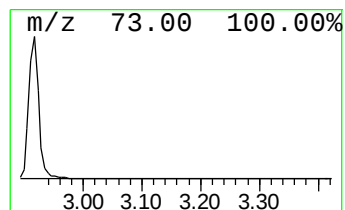
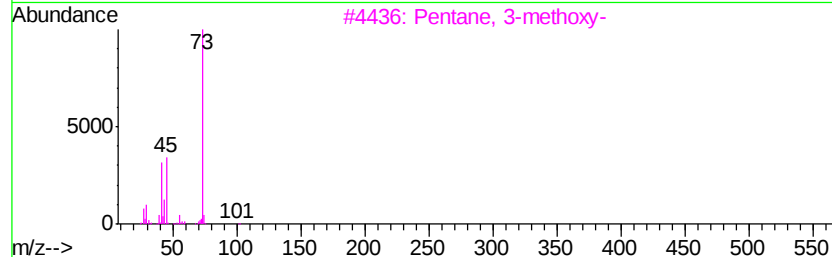
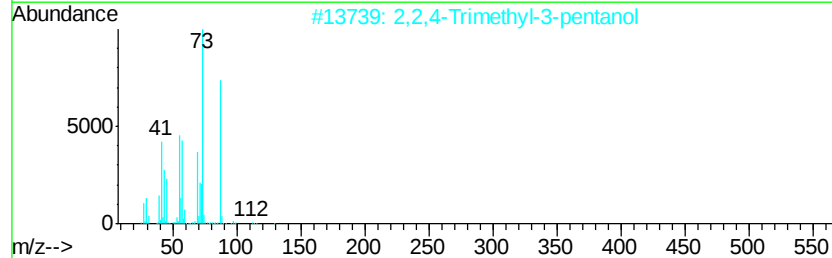
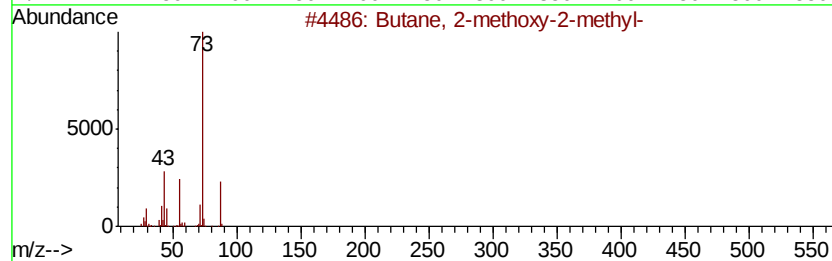
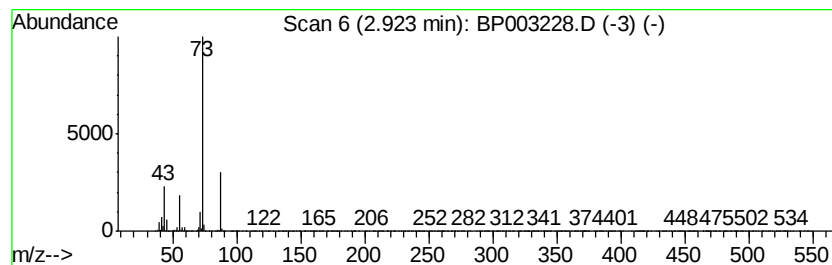
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	22.34 ng	1241960	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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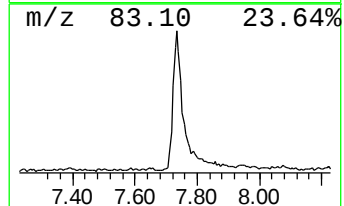
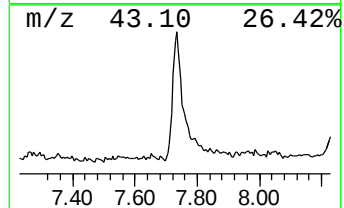
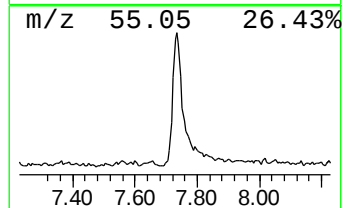
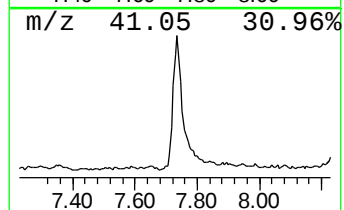
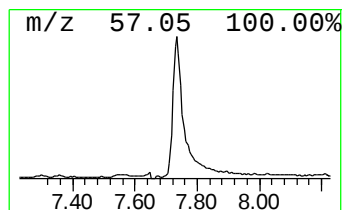
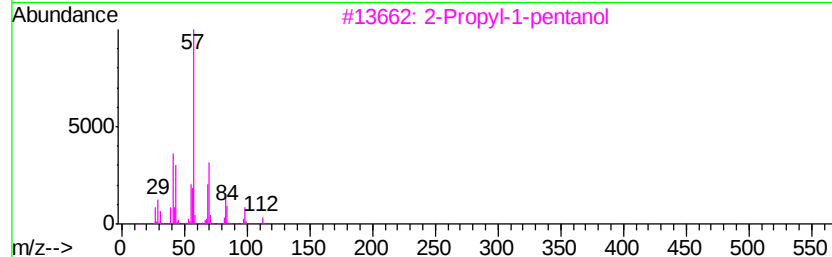
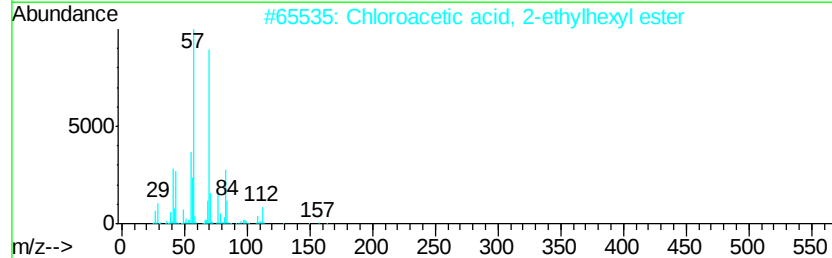
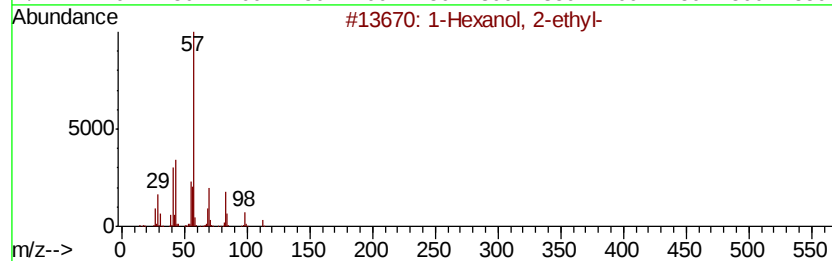
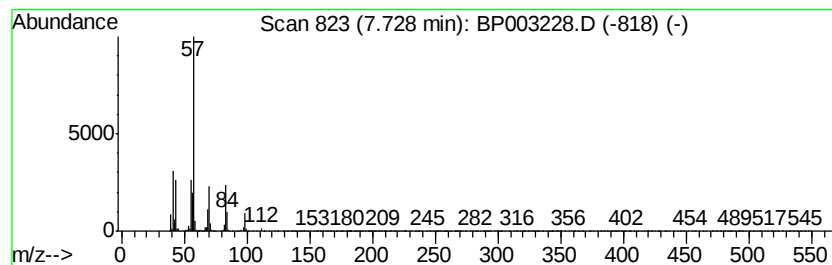
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	6.31 ng	350960	1,4-Dichlorobenzene-d4	7.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Chloroacetic acid, 2-ethylhexyl ...	206	C10H19ClO2	005345-58-4	59
3			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50
4			2-Decene, 5-methyl-, (Z)-	154	C11H22	074645-86-6	47
5			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	42



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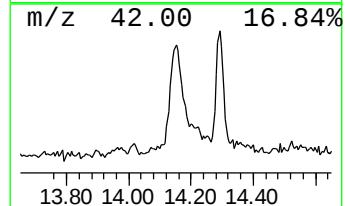
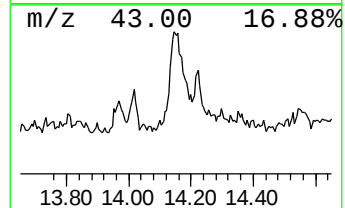
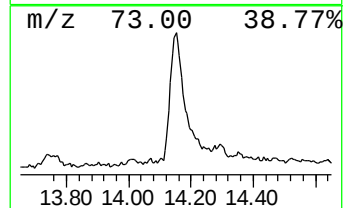
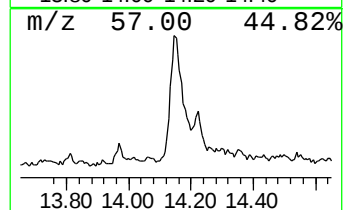
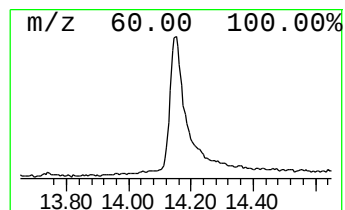
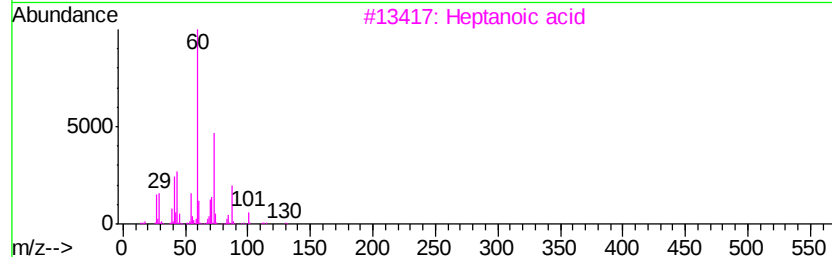
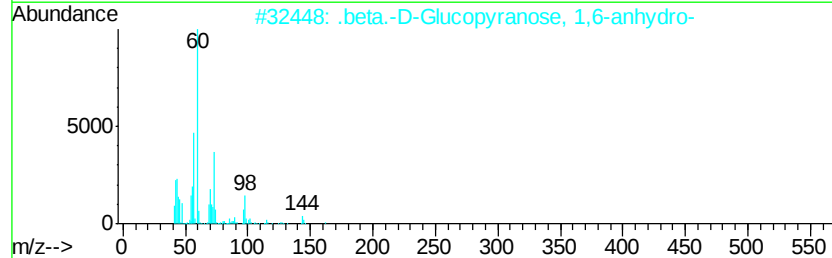
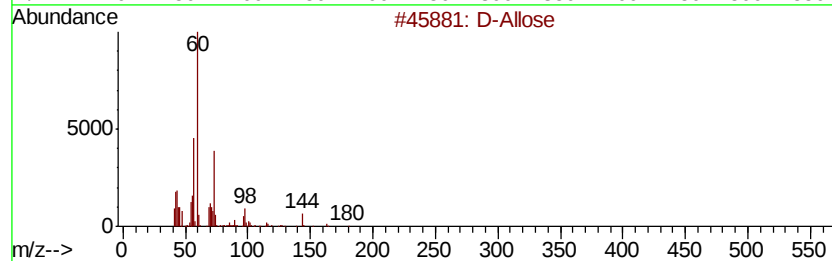
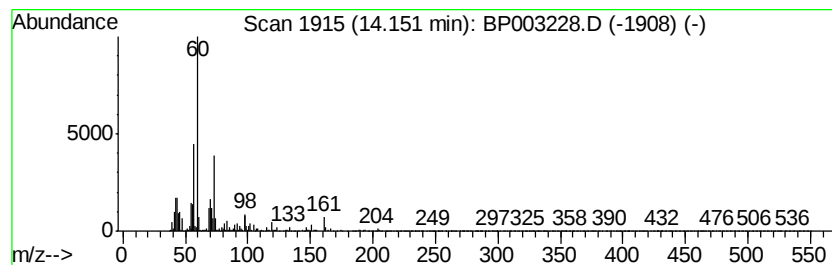
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 D-Allose Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	3.39 ng	279438	Acenaphthene-d10	14.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	D-Allose		180 C6H12O6	002595-97-3 80
2	.beta.-D-Glucopyranose, 1,6-anhy...		162 C6H10O5	000498-07-7 72
3	Heptanoic acid		130 C7H14O2	000111-14-8 64
4	Ethyl .beta.-d-ribose		178 C7H14O5	1000126-95-4 59
5	Hexanoic acid		116 C6H12O2	000142-62-1 59



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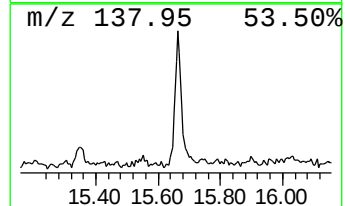
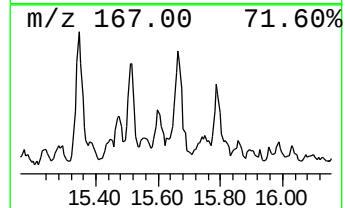
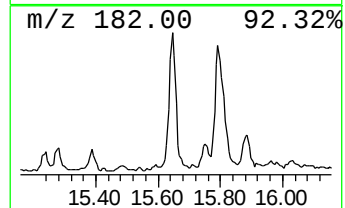
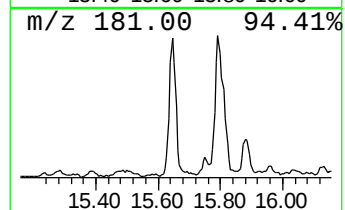
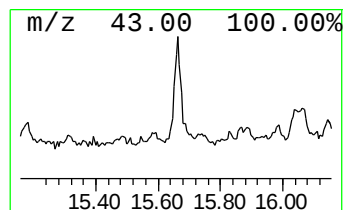
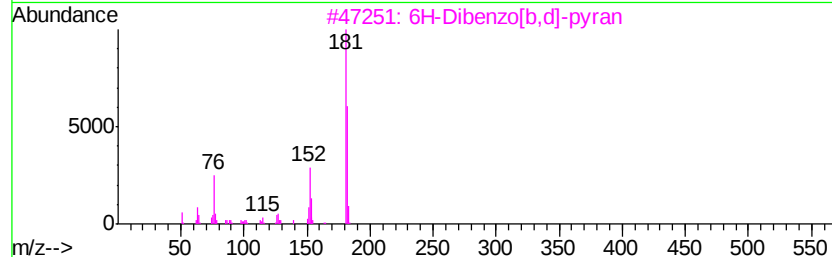
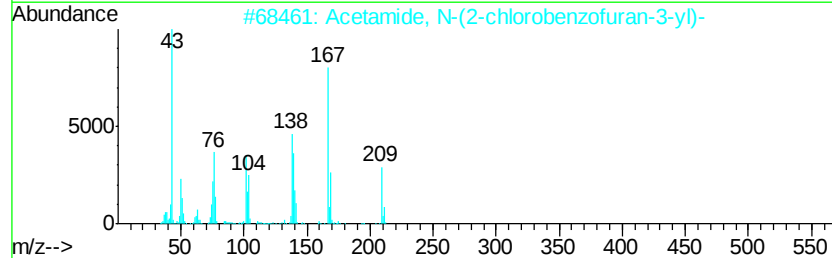
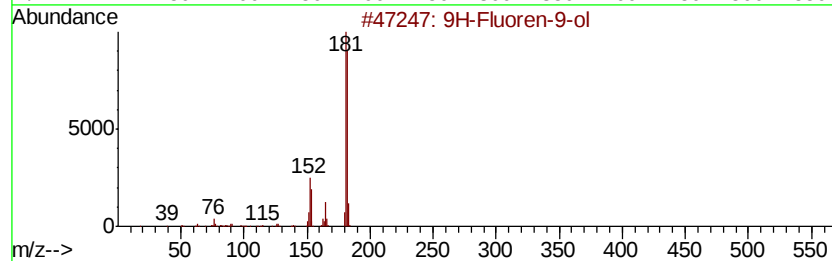
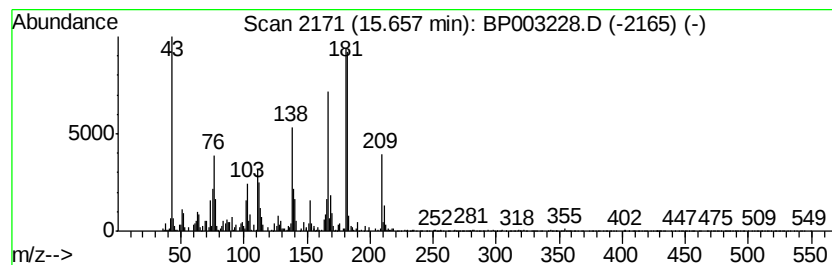
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9H-Fluoren-9-ol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	2.31 ng	190528	Acenaphthene-d10	14.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	9H-Fluoren-9-ol	182 C13H10O	001689-64-1	50
2	Acetamide, N-(2-chlorobenzofuran...	209 C10H8ClNO2	067382-11-0	22
3	6H-Dibenzo[b,d]-pyran	182 C13H10O	000229-95-8	22
4	Benzene, 1,2,3-trimethoxy-5-methyl-	182 C10H14O3	006443-69-2	22
5	Naphthalen-2-ol, 1-[3-(4-nitroph...	318 C19H14N2O3	1000294-84-3	20



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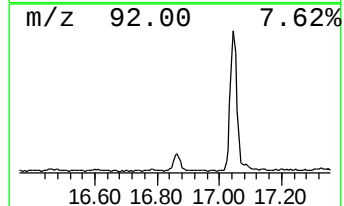
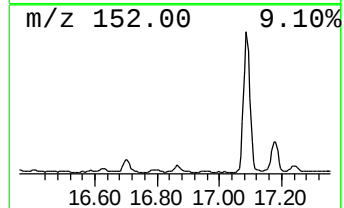
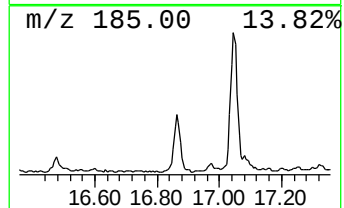
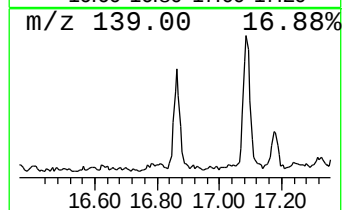
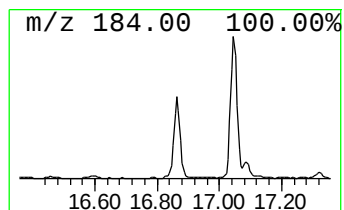
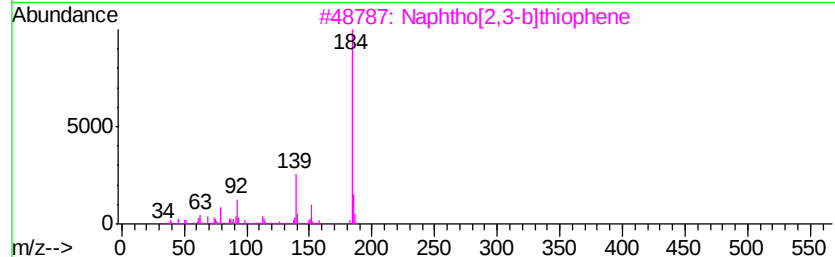
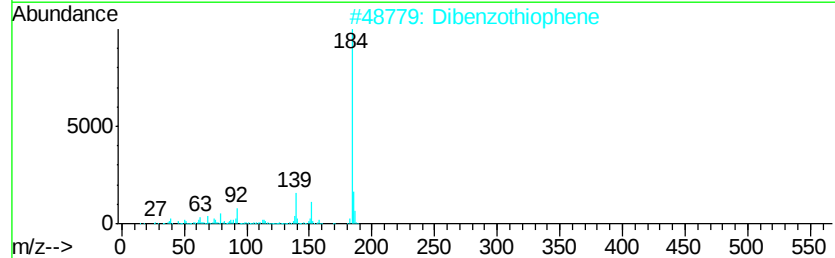
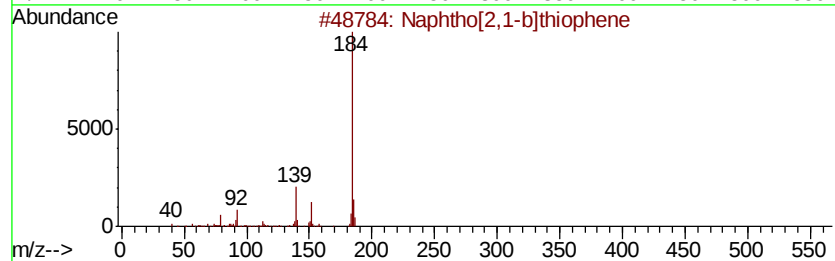
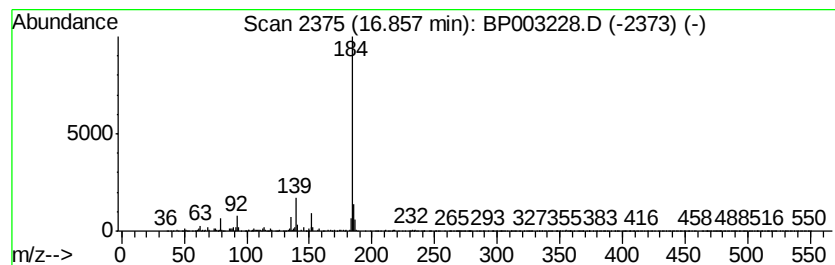
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphtho[2,1-b]thiophene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.86	2.19 ng	225573	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	96
2		Dibenzothiophene	184	C12H8S	000132-65-0	95
3		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	93
4		Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	90
5		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
Operator : CG/JU
Sample : L3916-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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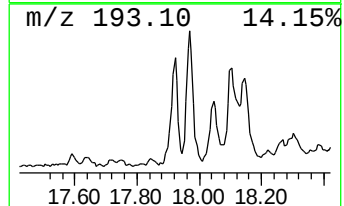
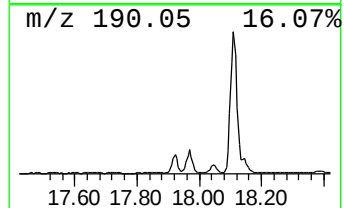
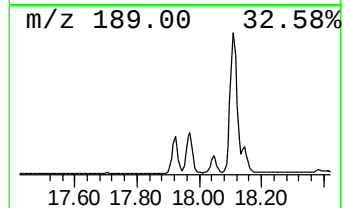
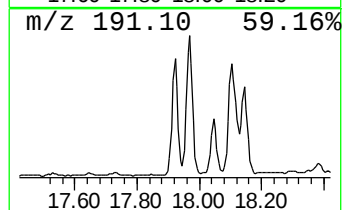
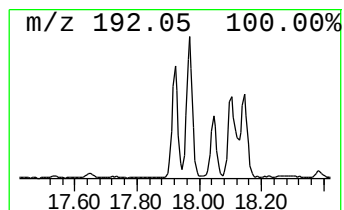
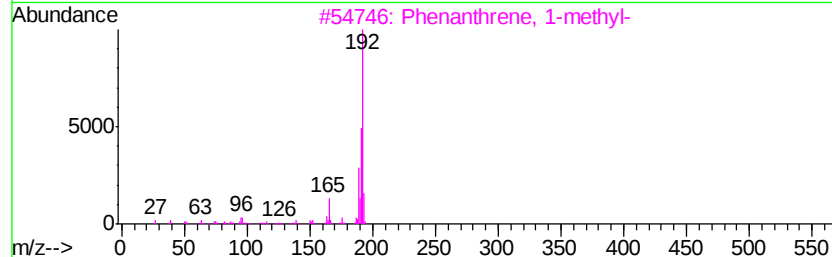
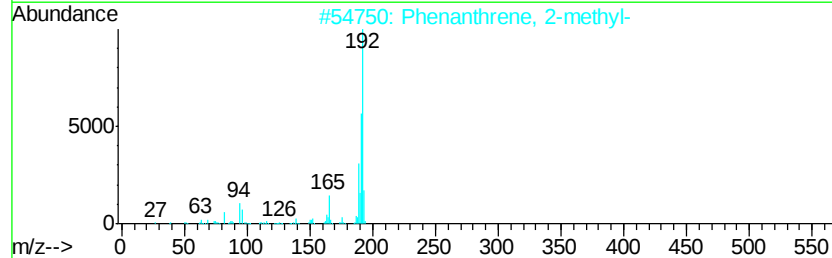
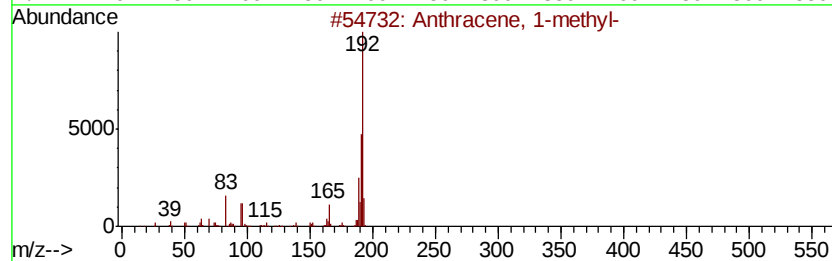
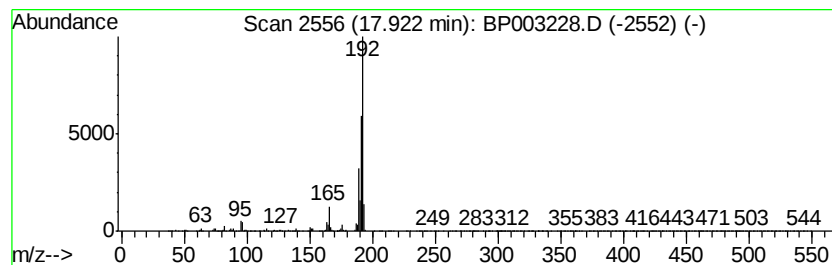
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	3.45 ng	355924	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
Operator : CG/JU
Sample : L3916-01
Misc :
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Instrument :
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ClientSampleId :
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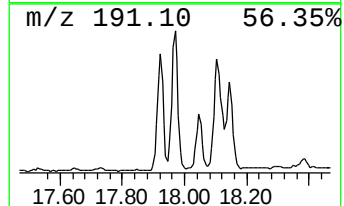
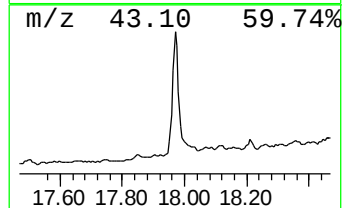
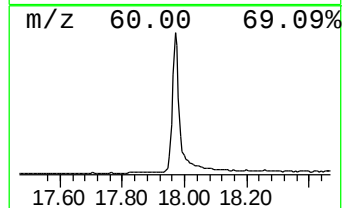
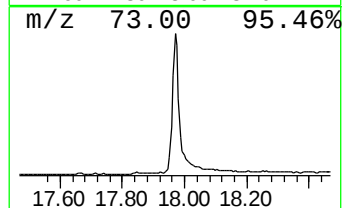
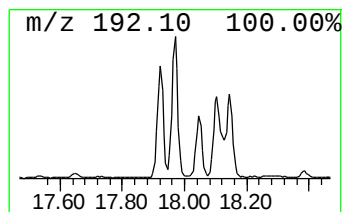
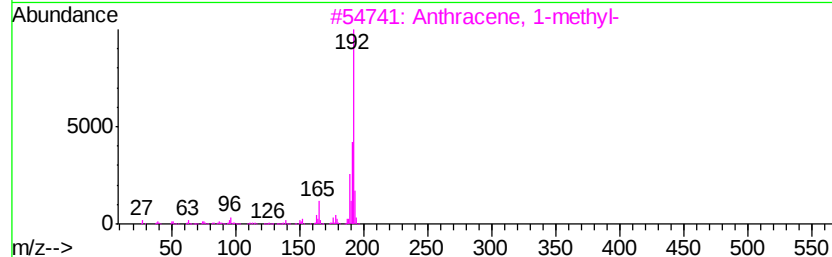
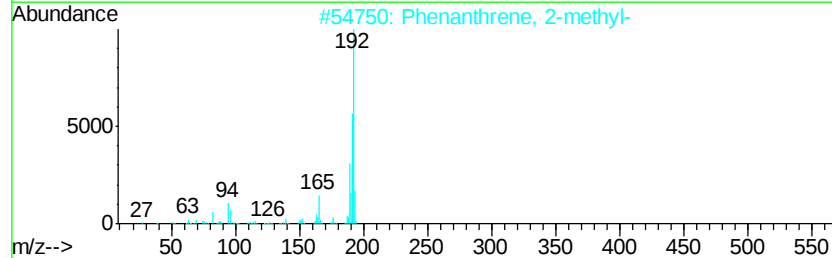
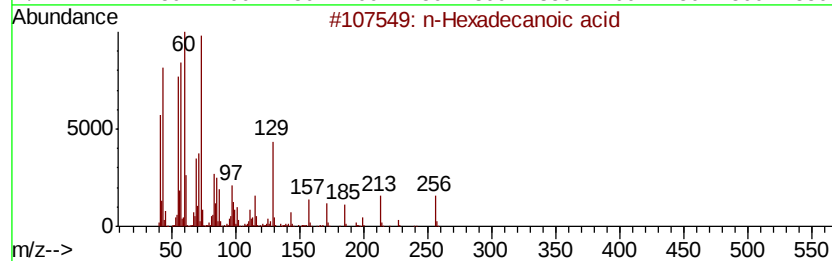
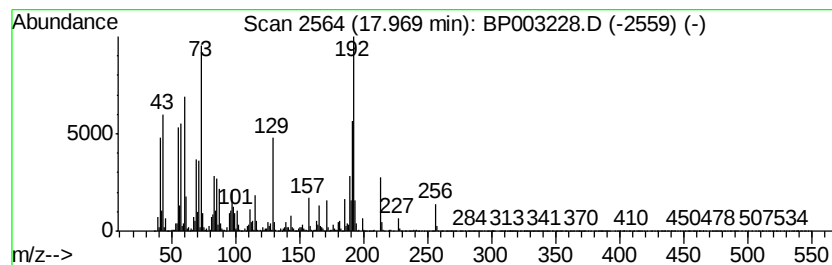
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	21.90 ng	2257620	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
Operator : CG/JU
Sample : L3916-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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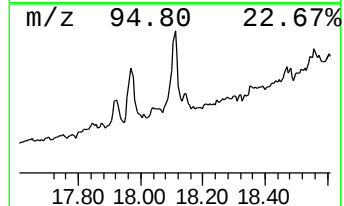
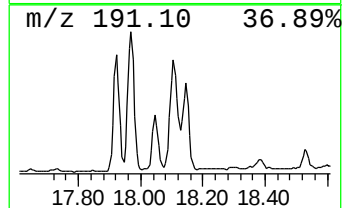
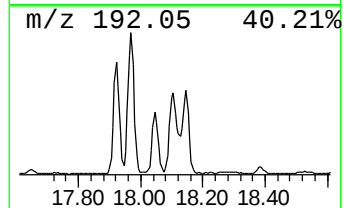
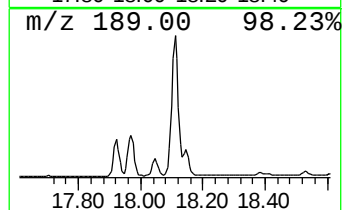
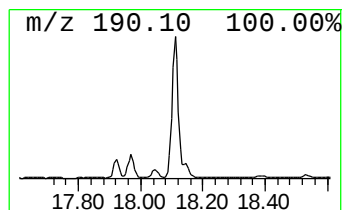
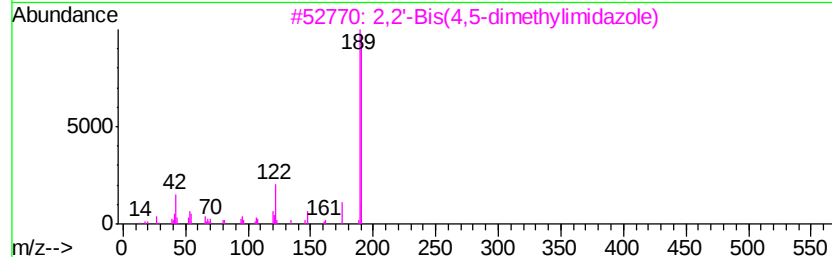
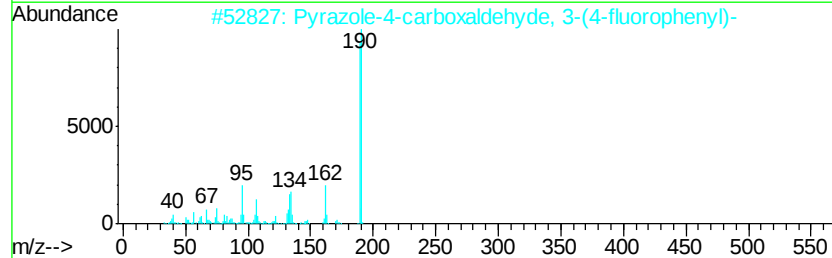
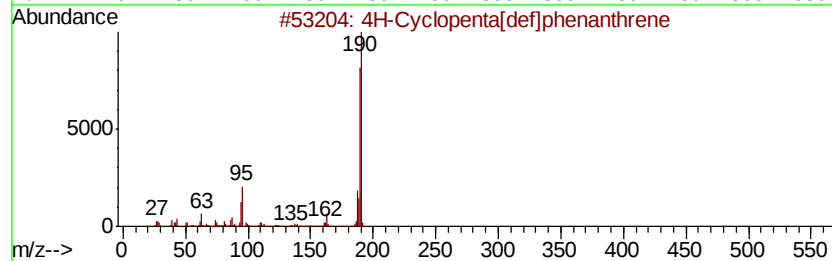
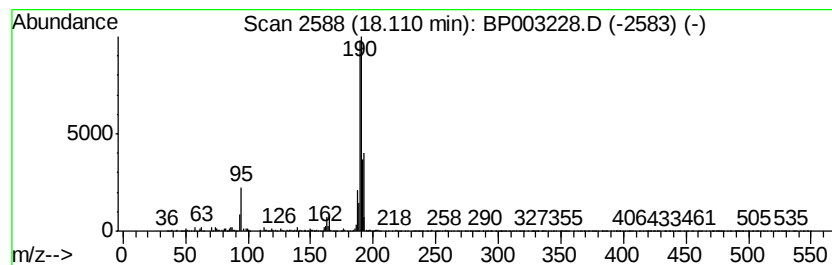
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	8.75 ng	902419	Phenanthrene-d10	17.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	86
2		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	52
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	49
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
5		2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	25



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Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
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Sample : L3916-01
Misc :
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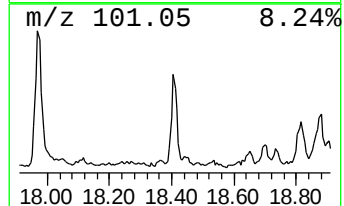
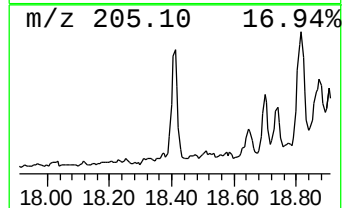
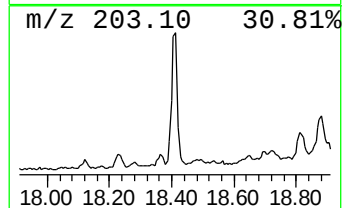
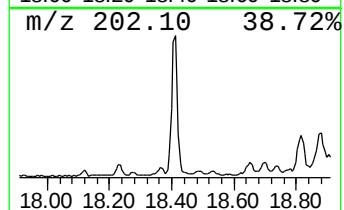
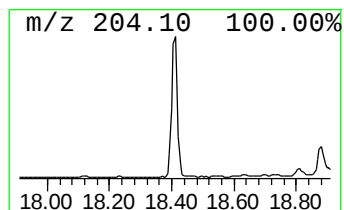
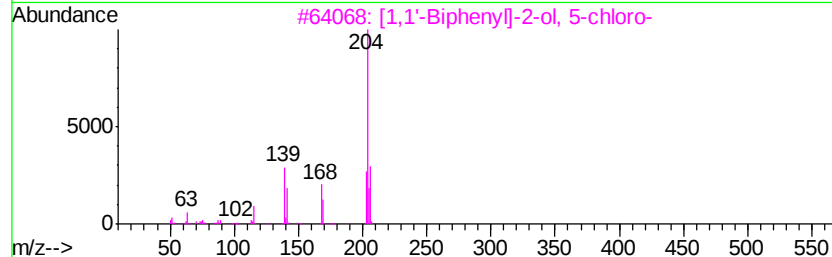
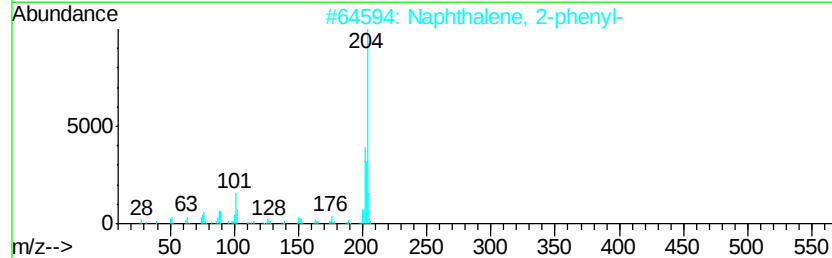
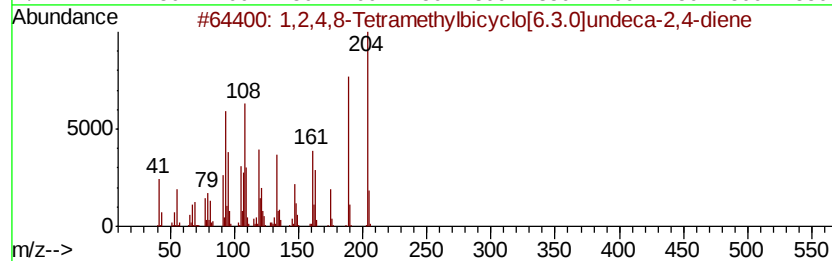
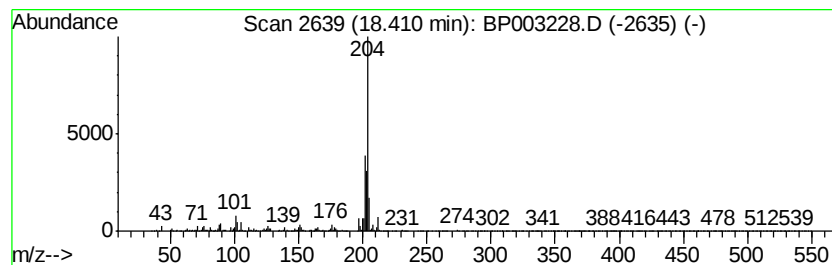
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1,2,4,8-Tetramethylbicyclo[... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	3.51 ng	362404	Phenanthrene-d10	17.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene	204	C15H24	137235-51-9	83
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
3			[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	47
4			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	46
5			5,16[1',2']:[8,13[1'',2'']]-Dibenz[1,2-b:4,5-b']diphenyl	408	C32H24	005672-97-9	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
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Sample : L3916-01
Misc :
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Instrument :
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ClientSampleId :
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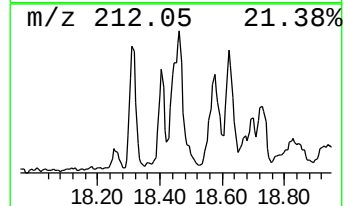
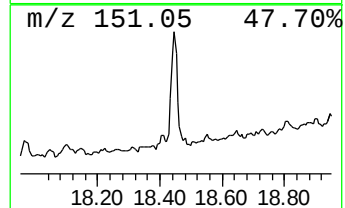
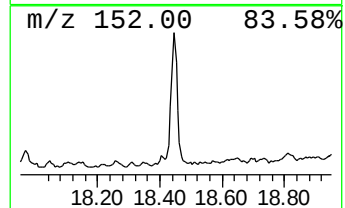
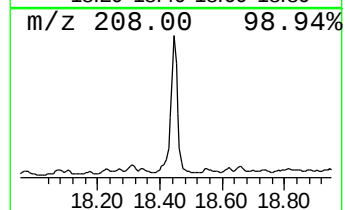
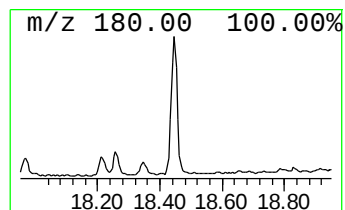
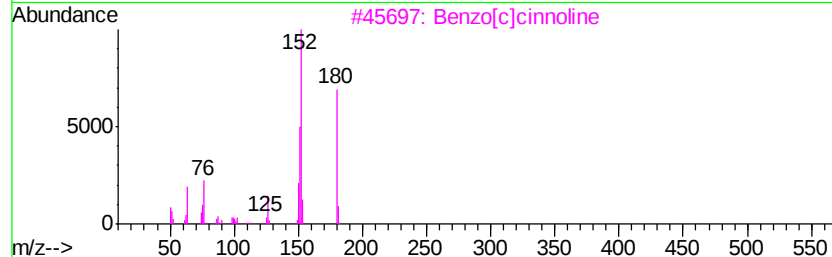
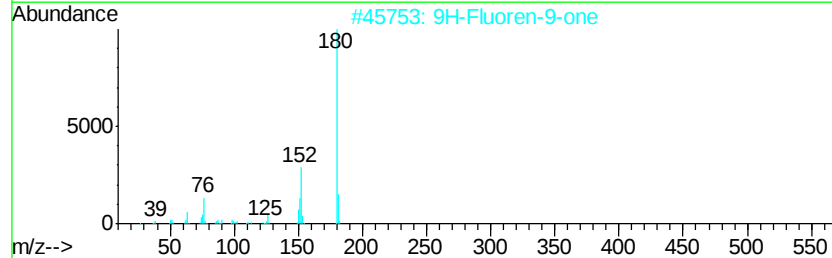
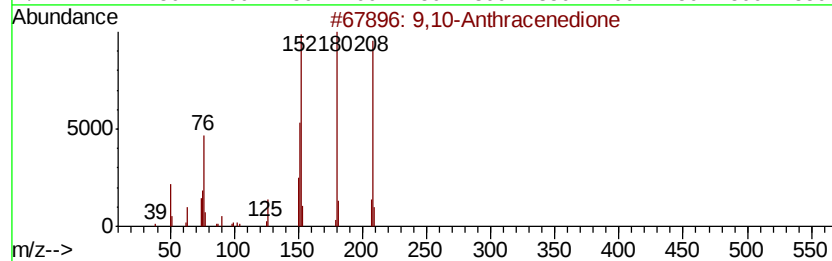
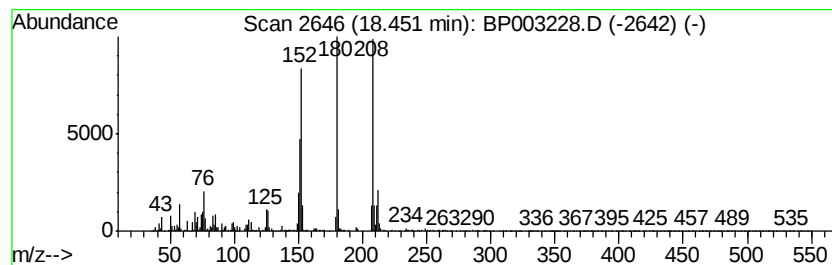
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.45	4.21 ng	433929	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2		9H-Fluoren-9-one	180	C13H8O	000486-25-9	83
3		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	64
4		1H-Phenalen-1-one	180	C13H8O	000548-39-0	60
5		1,4-Anthracenedione	208	C14H8O2	000635-12-1	52



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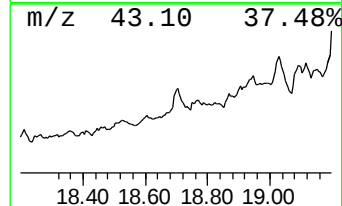
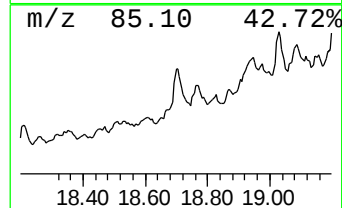
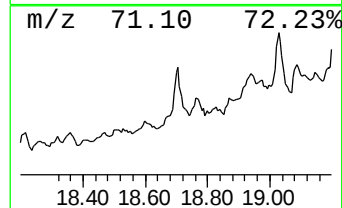
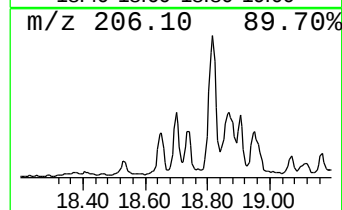
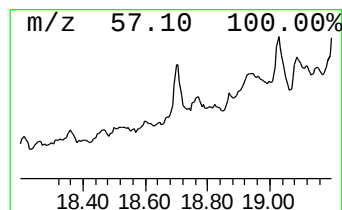
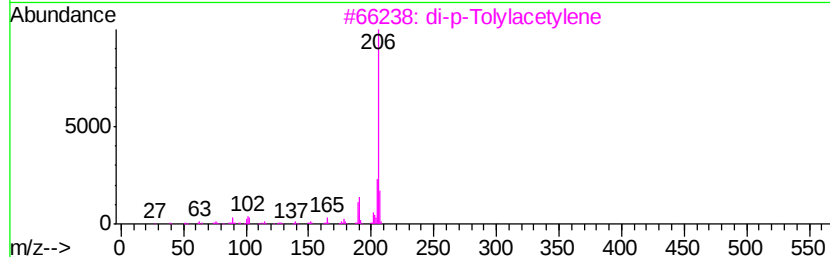
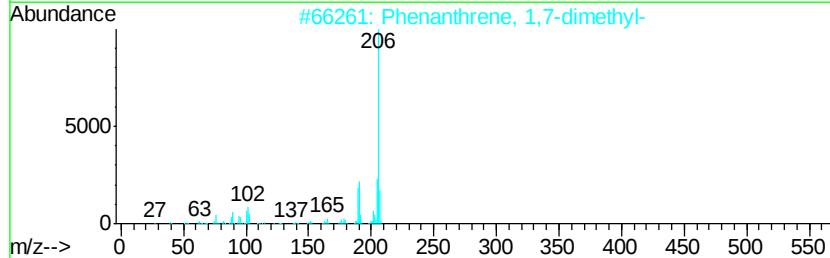
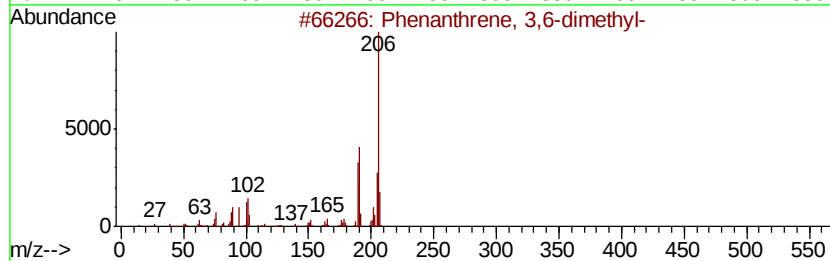
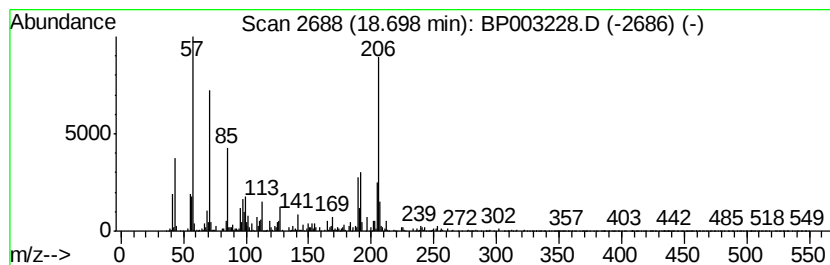
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenanthrene, 3,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.70	2.21 ng	227404	Phenanthrene-d10	17.05	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	92
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	83
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	64
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	60
5	Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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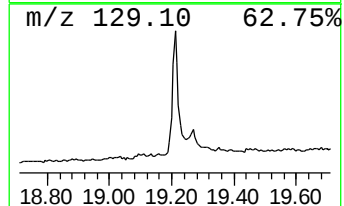
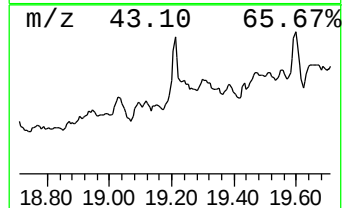
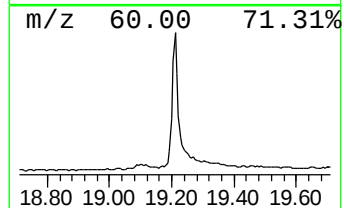
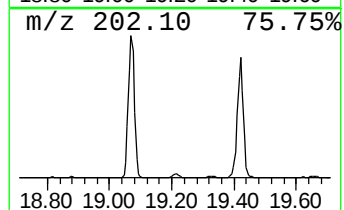
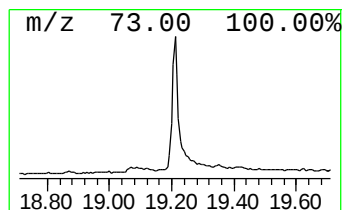
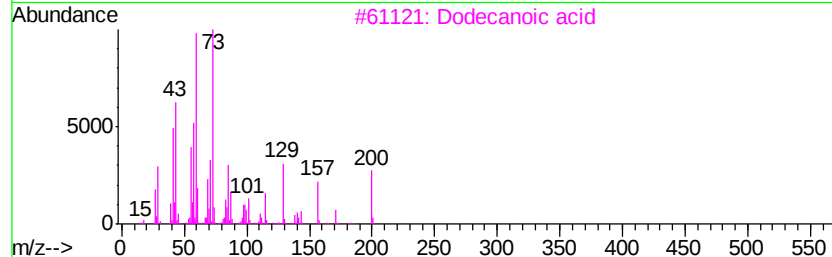
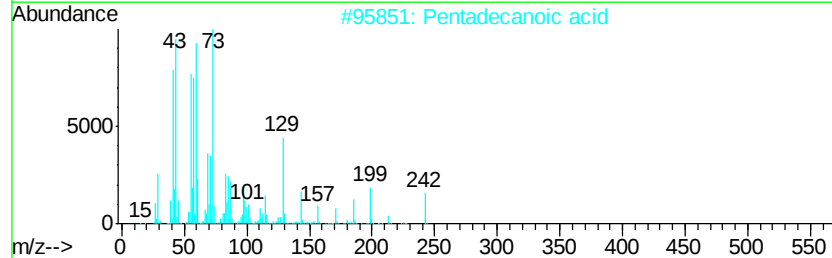
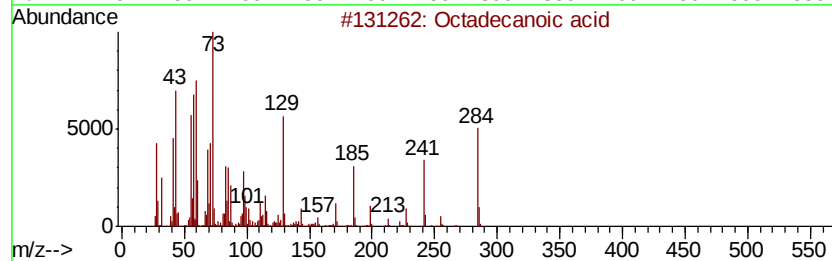
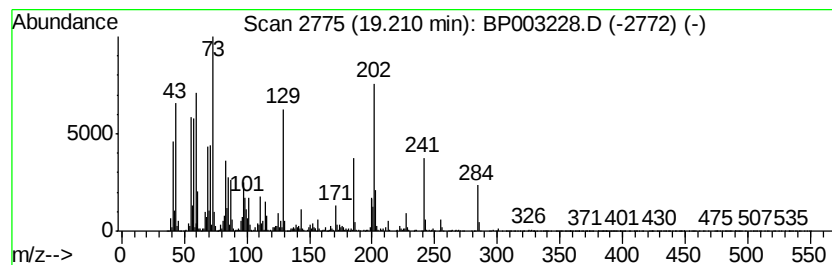
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.21	5.55 ng	1129400	Chrysene-d12	21.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3		Dodecanoic acid	200	C12H24O2	000143-07-7	47
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	45
5		Tridecanoic acid	214	C13H26O2	000638-53-9	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
Operator : CG/JU
Sample : L3916-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WC-1

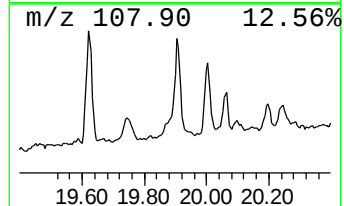
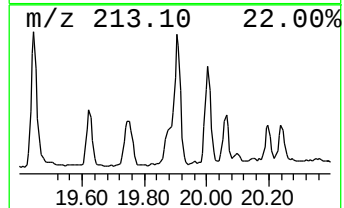
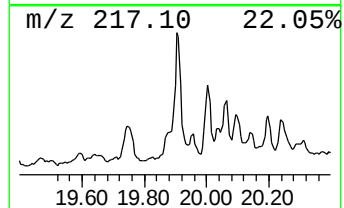
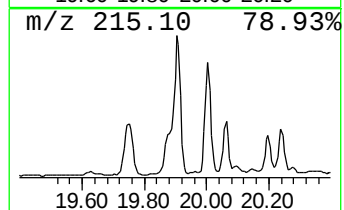
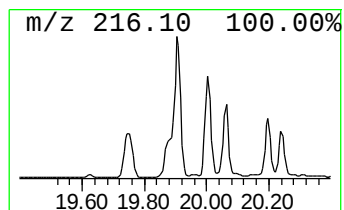
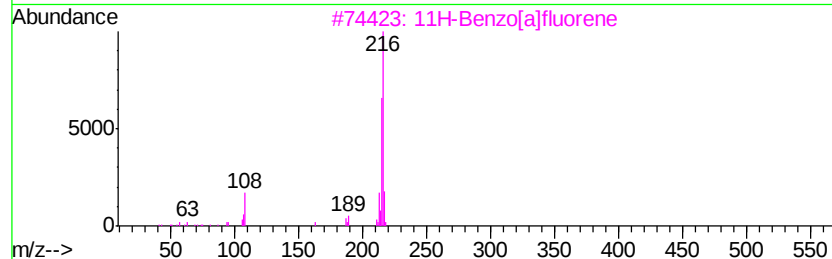
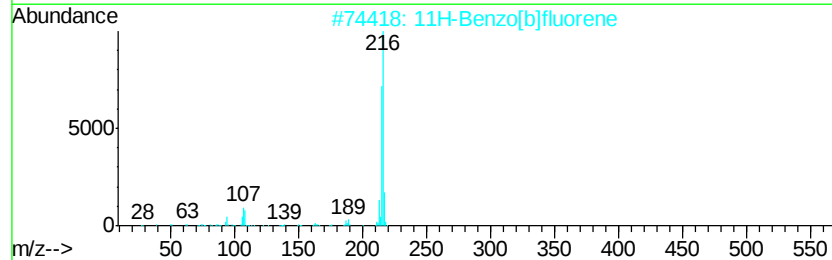
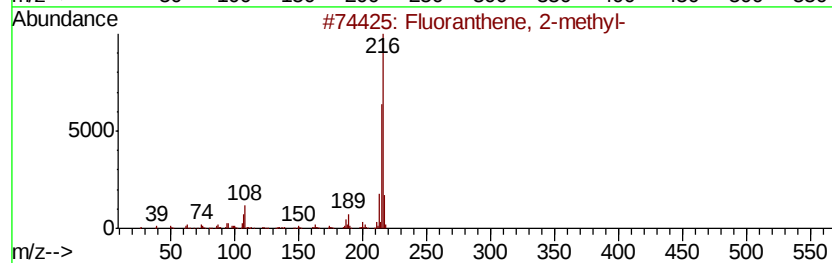
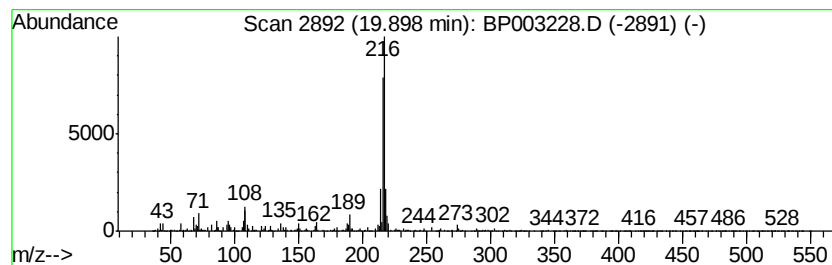
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Fluoranthene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.90	4.47 ng	909443	Chrysene-d12	21.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP090820\
Data File : BP003228.D
Acq On : 08 Sep 2020 15:48
Operator : CG/JU
Sample : L3916-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.92	22.3	ng	1241960	1	7.65	1111990	20.0
1-Hexanol, 2-ethyl-	7.73	6.3	ng	350960	1	7.65	1111990	20.0
D-Allose	14.15	3.4	ng	279438	3	14.29	1650130	20.0
9H-Fluoren-9-ol	15.66	2.3	ng	190528	3	14.29	1650130	20.0
Naphtho[2,1-b]thi...	16.86	2.2	ng	225573	4	17.05	2062110	20.0
Anthracene, 1-met...	17.92	3.5	ng	355924	4	17.05	2062110	20.0
n-Hexadecanoic acid	17.97	21.9	ng	2257620	4	17.05	2062110	20.0
4H-Cyclopenta[def...	18.11	8.8	ng	902419	4	17.05	2062110	20.0
1,2,4,8-Tetrameth...	18.41	3.5	ng	362404	4	17.05	2062110	20.0
9,10-Anthracenedione	18.45	4.2	ng	433929	4	17.05	2062110	20.0
Phenanthrene, 3,6...	18.70	2.2	ng	227404	4	17.05	2062110	20.0
Octadecanoic acid	19.21	5.5	ng	1129400	5	21.15	4068570	20.0
Fluoranthene, 2-m...	19.90	4.5	ng	909443	5	21.15	4068570	20.0