

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP005723.D
 Data File : BP005723.D
 Acq On : 04 Jun 2021 21:42
 Operator : CG/JU
 Sample : M2589-13 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B10(4-6)

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-BP052721.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005723.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.464	230	234	240	rBV	59861	81867	1.21%	0.162%
2	5.075	332	338	346	rBV	280445	394155	5.81%	0.781%
3	5.546	412	418	425	rBV	431544	647111	9.54%	1.282%
4	6.663	602	608	615	rVB	87998	142021	2.09%	0.281%
5	7.152	681	691	706	rBV	845740	1460401	21.52%	2.894%
6	7.987	824	833	839	rBV	572535	950816	14.01%	1.884%
7	9.146	1024	1030	1038	rVV	590882	1012243	14.92%	2.006%
8	9.216	1038	1042	1053	rVB3	49045	95427	1.41%	0.189%
9	10.793	1300	1310	1315	rBV	751949	1349914	19.89%	2.675%
10	10.846	1315	1319	1326	rVB	227466	384004	5.66%	0.761%
11	12.210	1542	1551	1557	rBV	65985	111750	1.65%	0.221%
12	12.446	1582	1591	1598	rVB	150815	258407	3.81%	0.512%
13	12.663	1621	1628	1638	rBV	114989	202330	2.98%	0.401%
14	13.240	1719	1726	1736	rVB	1567523	2354713	34.70%	4.667%
15	13.440	1753	1760	1767	rVB2	118457	199542	2.94%	0.395%
16	13.775	1812	1817	1824	rBV2	60589	158043	2.33%	0.313%
17	13.934	1839	1844	1849	rBV	113549	199078	2.93%	0.395%
18	13.987	1849	1853	1858	rVB	72448	109527	1.61%	0.217%
19	14.092	1868	1871	1875	rVB	56690	68256	1.01%	0.135%
20	14.334	1904	1912	1921	rBV	134417	243778	3.59%	0.483%
21	14.504	1937	1941	1947	rVB	129155	158438	2.33%	0.314%
22	14.610	1950	1959	1965	rBV	1171203	1735357	25.57%	3.439%
23	14.675	1965	1970	1978	rVB	363116	554677	8.17%	1.099%
24	14.822	1989	1995	2002	rVB	36488	97092	1.43%	0.192%
25	15.004	2021	2026	2032	rVB	234437	331678	4.89%	0.657%
26	15.081	2036	2039	2043	rVB	55291	70891	1.04%	0.140%
27	15.222	2060	2063	2068	rVB	46858	71459	1.05%	0.142%
28	15.392	2086	2092	2096	rBV3	52797	102468	1.51%	0.203%
29	15.451	2099	2102	2107	rVB2	116745	138045	2.03%	0.274%
30	15.651	2131	2136	2148	rBV	333143	571756	8.43%	1.133%
31	15.816	2161	2164	2168	rBV3	57785	75341	1.11%	0.149%
32	15.857	2168	2171	2175	rVB2	70859	97596	1.44%	0.193%
33	15.945	2183	2186	2194	rVB3	85010	157632	2.32%	0.312%
34	16.092	2207	2211	2218	rVB2	184869	312503	4.61%	0.619%
35	16.316	2245	2249	2250	rBV	157729	197192	2.91%	0.391%
36	16.651	2300	2306	2311	rBV4	114387	215179	3.17%	0.426%

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

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37	17.010	2363	2367	2373	rVB2	73346	104433	1.54%	0.207%
38	17.104	2380	2383	2388	rVB	152123	182085	2.68%	0.361%
39	17.169	2390	2394	2404	rVB2	212153	356050	5.25%	0.706%
40	17.357	2420	2426	2429	rBV	1345045	2020508	29.78%	4.004%
41	17.398	2429	2433	2438	rVB	2739935	3782557	55.74%	7.497%
42	17.486	2444	2448	2454	rVB	766544	1027001	15.13%	2.035%
43	17.545	2455	2458	2462	rVV2	63285	93138	1.37%	0.185%
44	17.757	2490	2494	2500	rVB	236388	328309	4.84%	0.651%
45	17.845	2506	2509	2512	rBV	145105	184940	2.73%	0.367%
46	18.216	2568	2572	2576	rBV	384724	517272	7.62%	1.025%
47	18.263	2576	2580	2585	rVB	515326	670269	9.88%	1.328%
48	18.339	2590	2593	2597	rVB	174405	209668	3.09%	0.416%
49	18.404	2599	2604	2608	rBV2	547527	969549	14.29%	1.922%
50	18.510	2619	2622	2625	rBV2	171450	200815	2.96%	0.398%
51	18.692	2650	2653	2657	rVB	232070	281954	4.16%	0.559%
52	19.357	2760	2766	2771	rBV	3602030	4884085	71.98%	9.680%
53	19.675	2816	2820	2822	rBV2	790071	1151922	16.98%	2.283%
54	19.710	2822	2826	2833	rVB	4891746	6785761	100.00%	13.449%
55	19.892	2853	2857	2862	rVB	2697165	3401711	50.13%	6.742%
56	20.186	2904	2907	2913	rVB2	644663	801194	11.81%	1.588%
57	21.416	3111	3116	3121	rBV3	2356731	4986658	73.49%	9.883%
58	21.463	3121	3124	3133	rVB	1671444	2206498	32.52%	4.373%

Sum of corrected areas: 50457064

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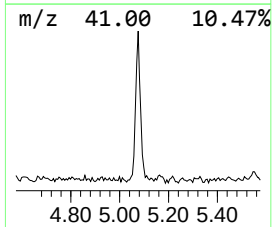
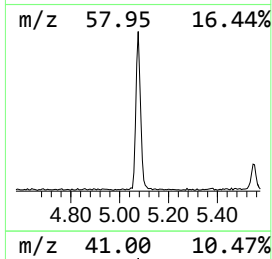
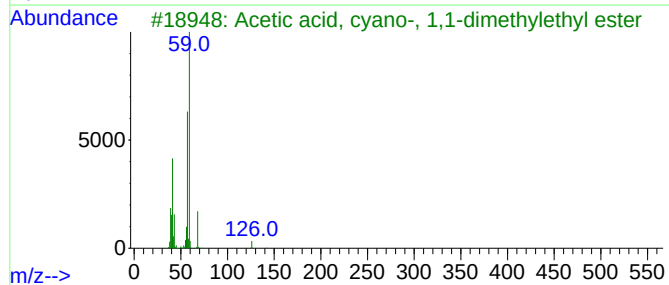
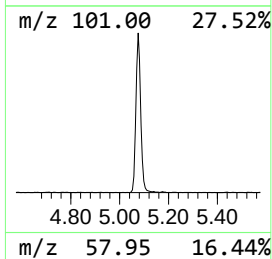
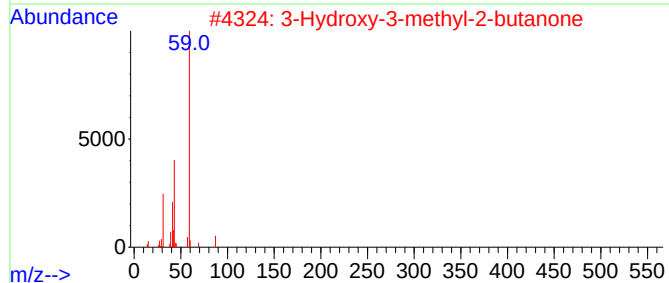
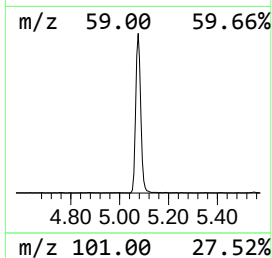
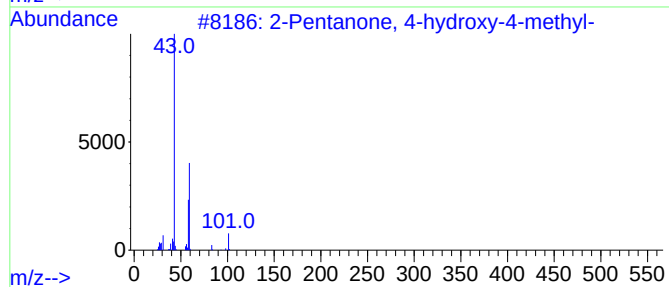
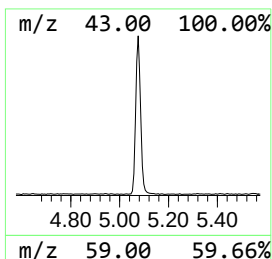
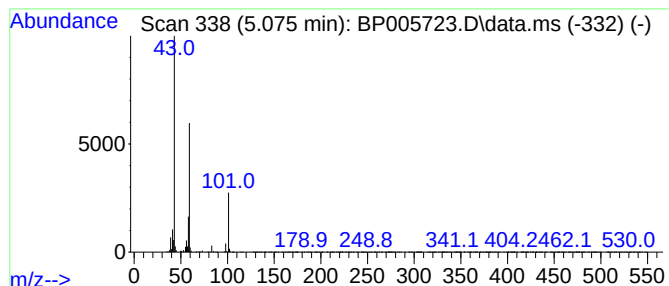
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	8.29 ng	394155	1,4-Dichlorobenzene-d4	7.987

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
4	2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	17		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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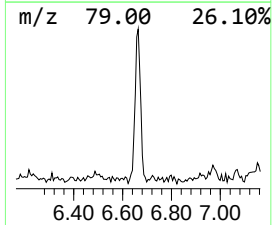
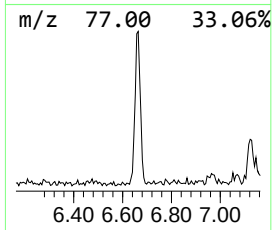
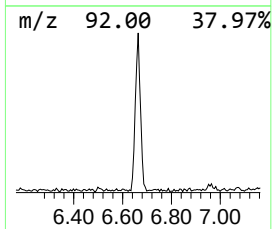
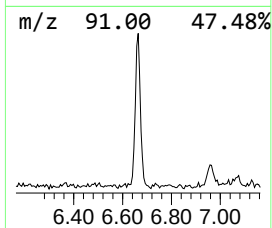
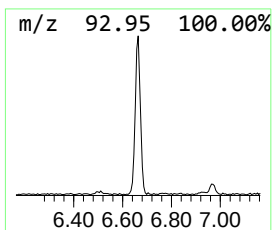
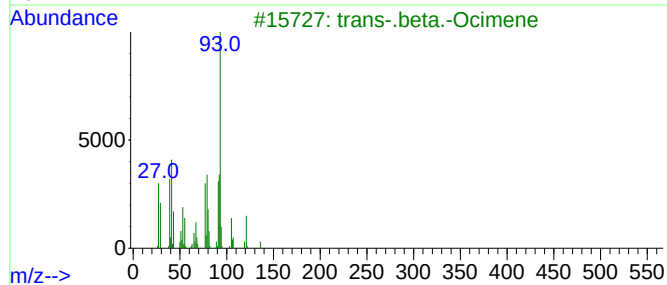
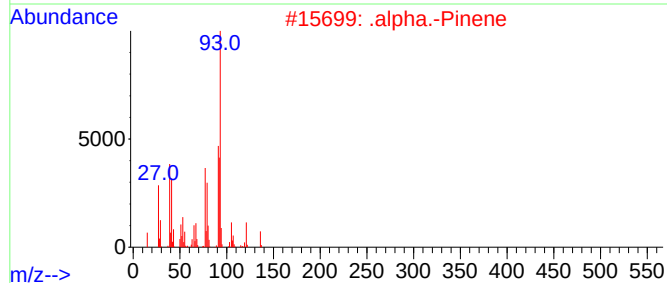
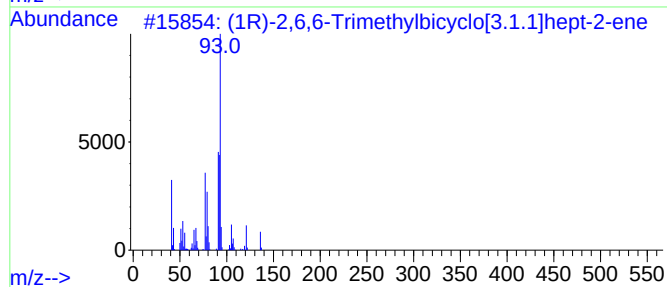
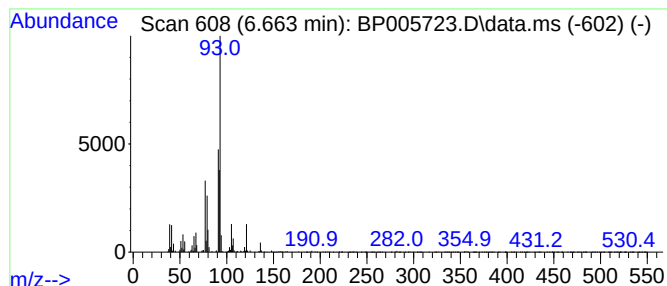
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.663	2.99 ng	142021	1,4-Dichlorobenzene-d4	7.987

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95		
2	.alpha.-Pinene	136	C10H16	000080-56-8	94		
3	trans-.beta.-Ocimene	136	C10H16	003779-61-1	94		
4	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91		
5	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	91		



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Instrument :
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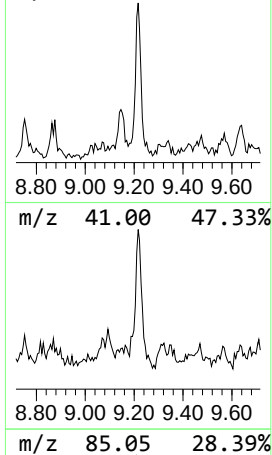
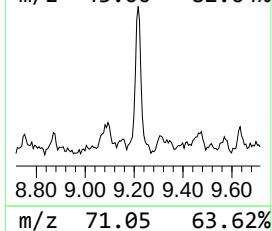
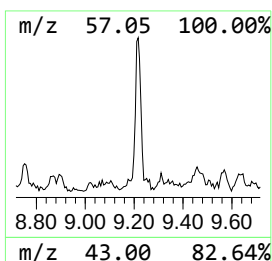
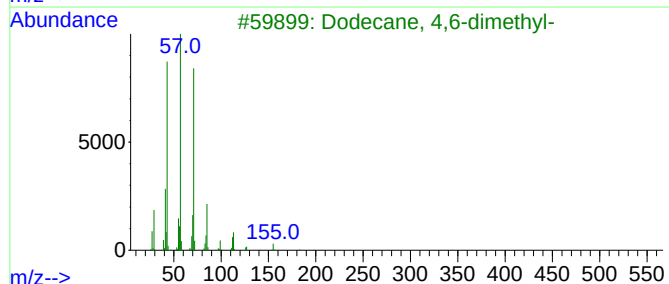
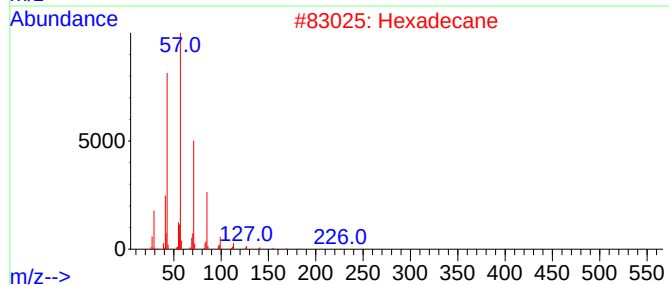
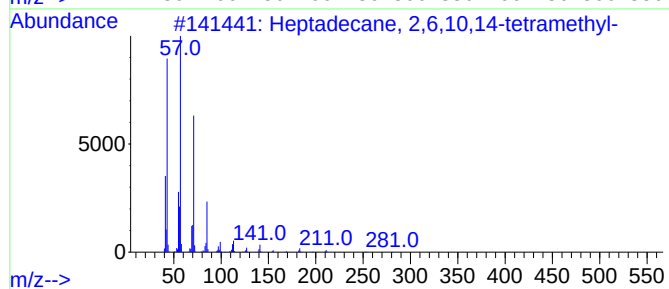
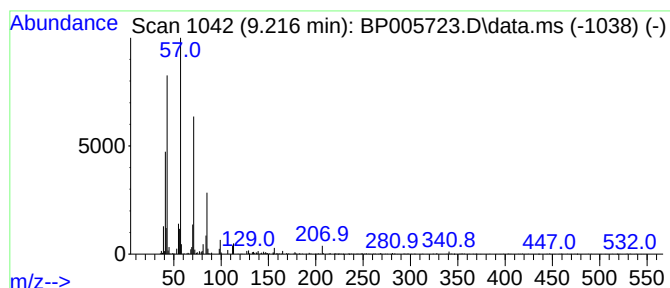
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Heptadecane, 2,6,10,14-tetr... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.216	2.01 ng	95427	1,4-Dichlorobenzene-d4	7.987

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	87
2			Hexadecane	226	C16H34	000544-76-3	72
3			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	68
4			Bacchotricuneatin c	342	C20H22O5	066563-30-2	68
5			Dodecane, 2-methyl-	184	C13H28	001560-97-0	59



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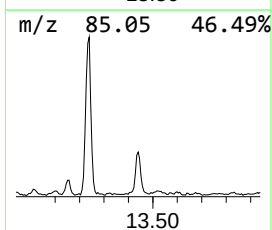
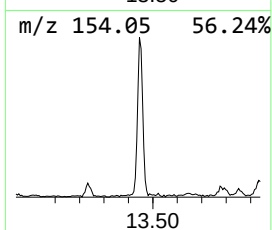
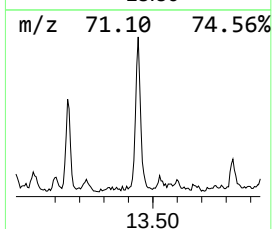
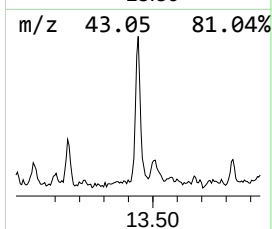
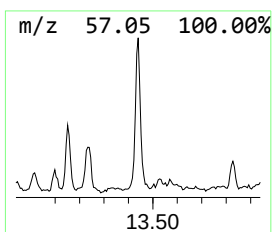
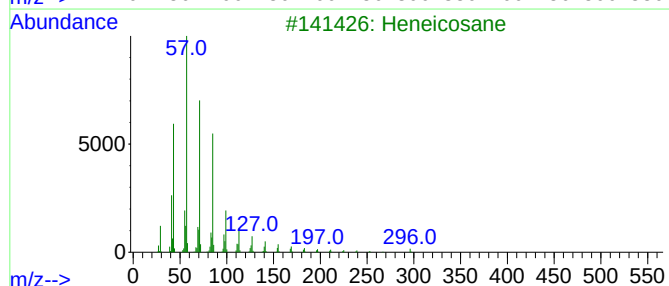
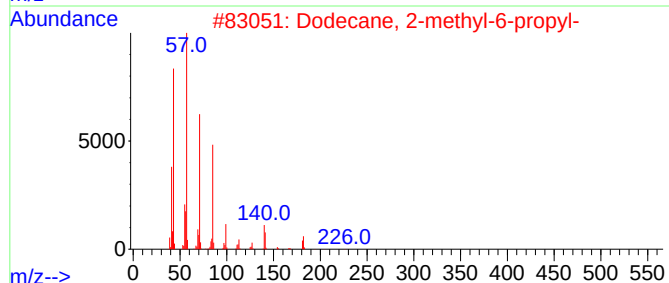
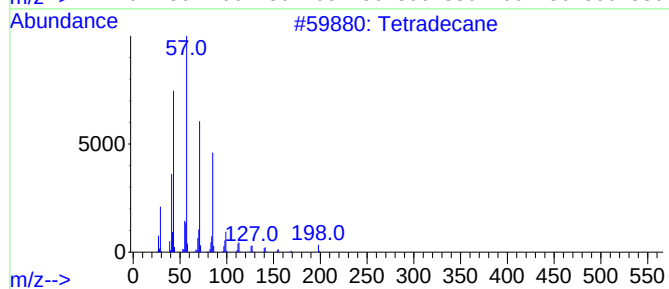
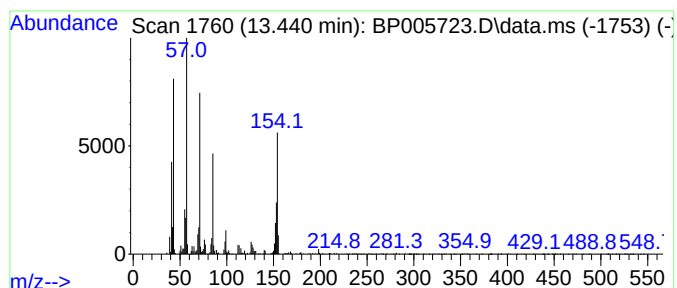
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.440	2.30 ng	199542	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	83
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	64
3			Heneicosane	296	C21H44	000629-94-7	60
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	58
5			Tridecane	184	C13H28	000629-50-5	58



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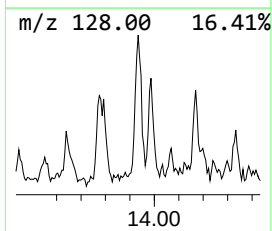
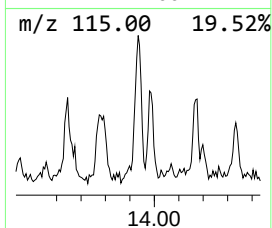
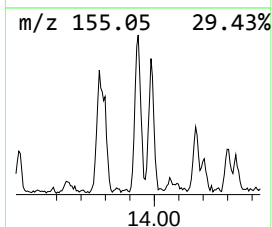
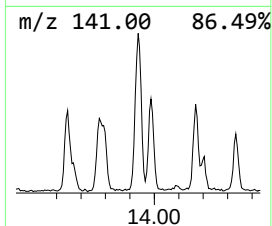
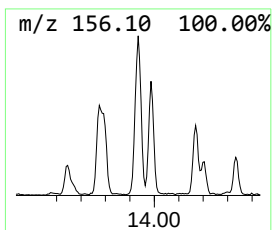
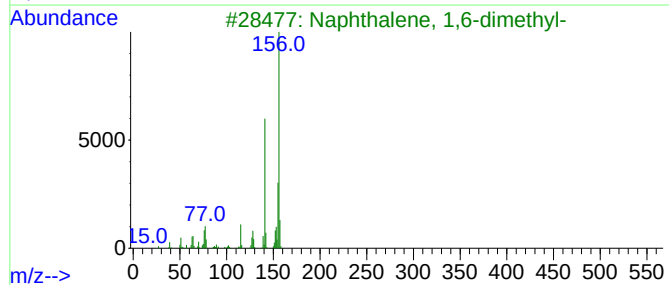
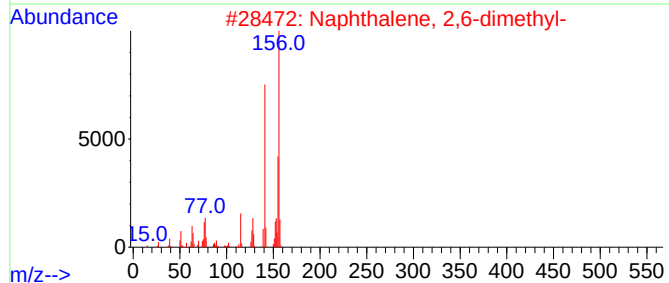
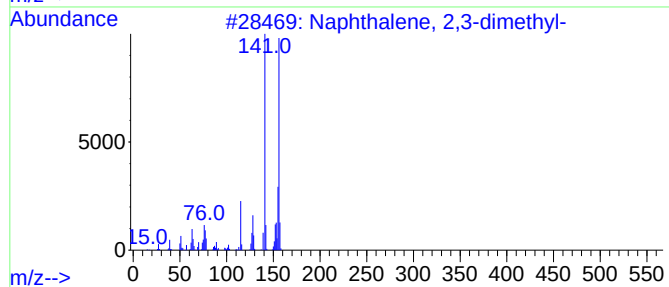
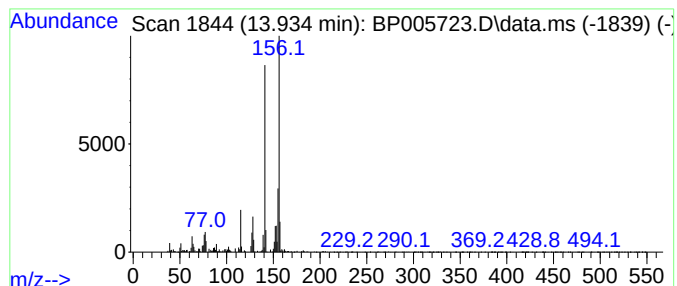
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.934	2.29 ng	199078	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
Sample : M2589-13 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B10(4-6)

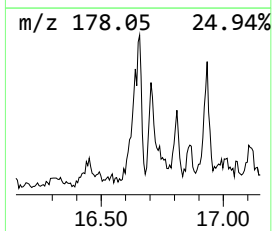
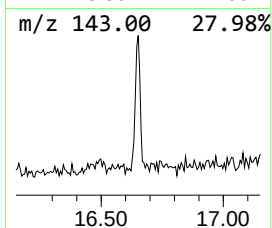
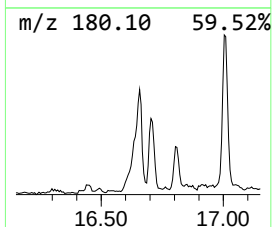
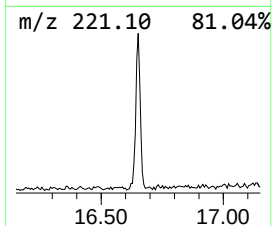
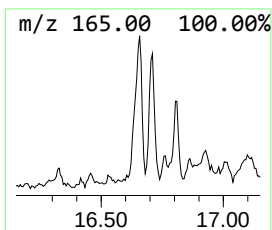
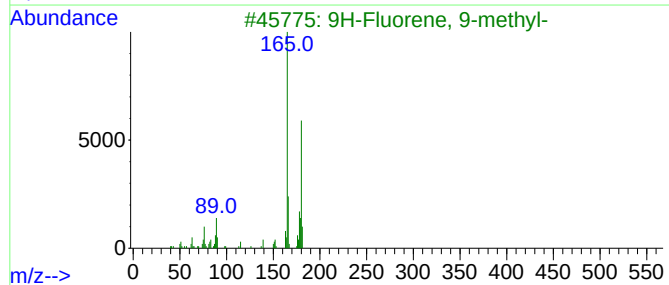
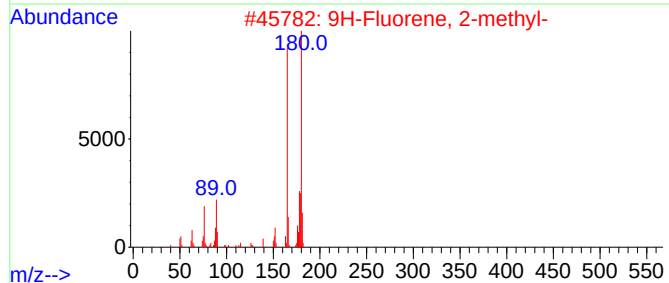
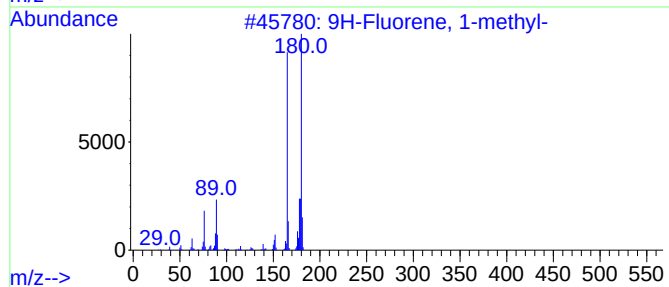
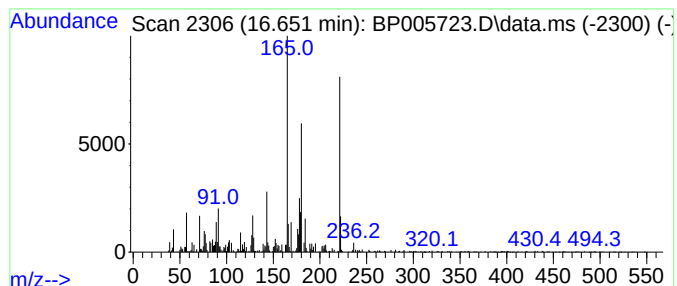
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9H-Fluorene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	2.13 ng	215179	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 1-methyl-			180	C14H12	001730-37-6	60
2	9H-Fluorene, 2-methyl-			180	C14H12	001430-97-3	60
3	9H-Fluorene, 9-methyl-			180	C14H12	002523-37-7	46
4	1H-Indene, 2,3-dihydro-1,1,3-tri...			236	C18H20	003910-35-8	41
5	3-Chloro-benzalacetone			180	C10H9ClO	1000146-90-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
Sample : M2589-13 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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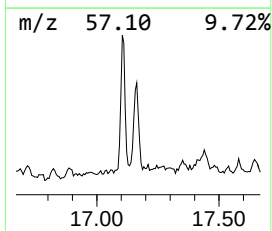
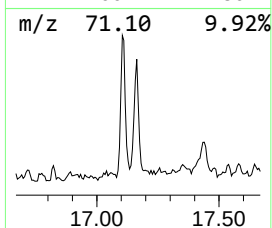
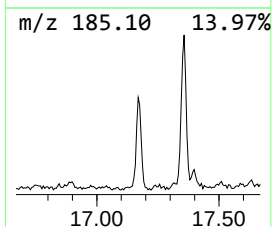
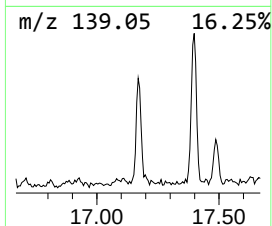
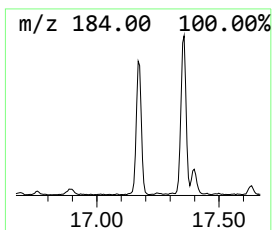
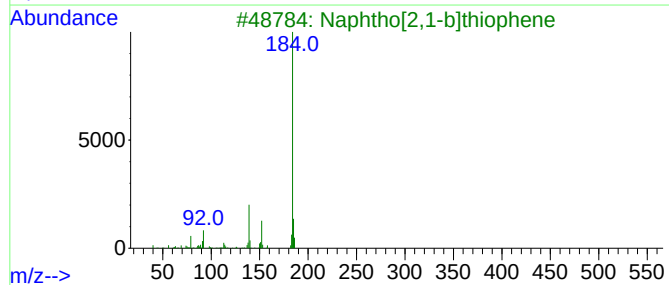
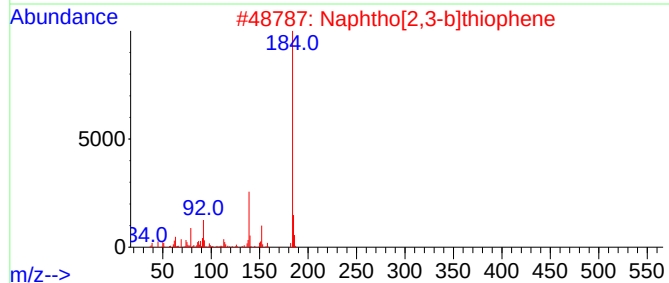
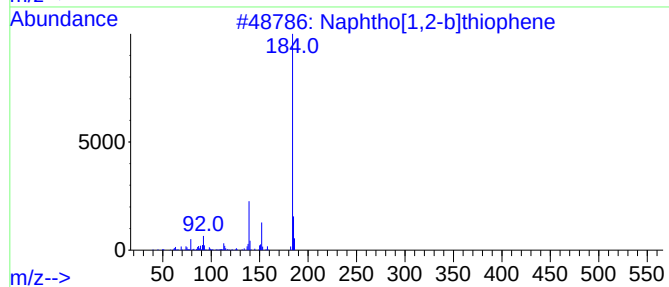
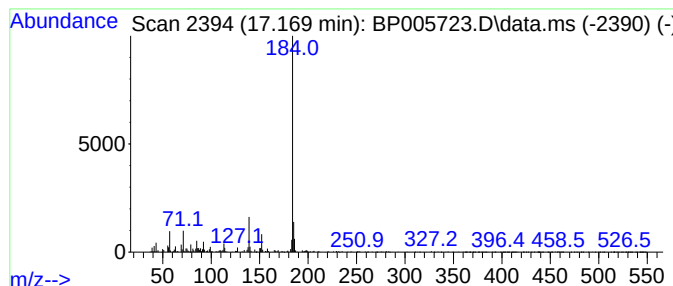
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphtho[1,2-b]thiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.169	3.52 ng	356050	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	94
2			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	93
3			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	93
4			Dibenzothiophene	184	C12H8S	000132-65-0	93
5			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
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ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
18-B10(4-6)

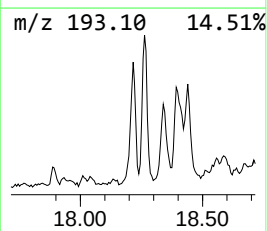
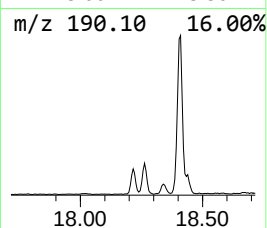
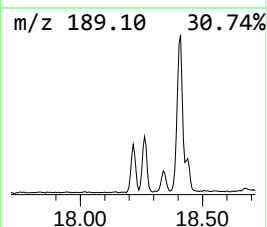
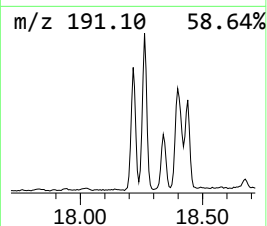
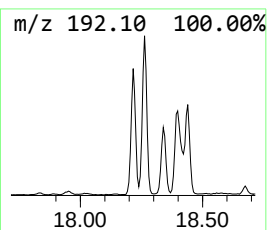
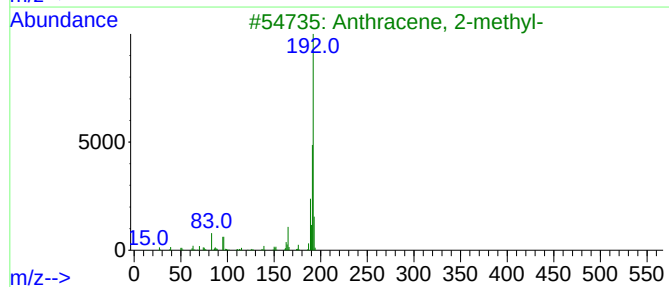
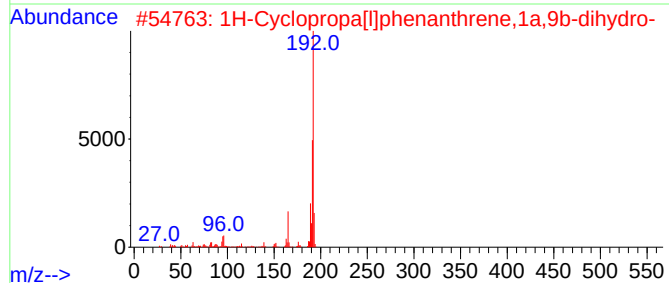
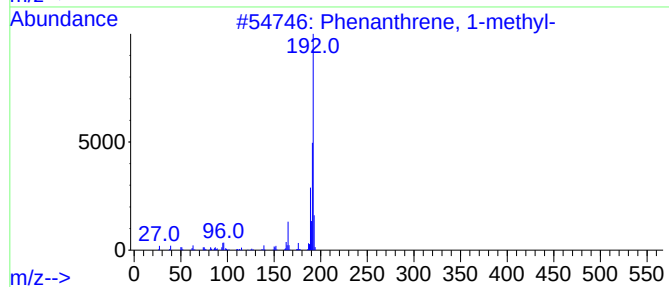
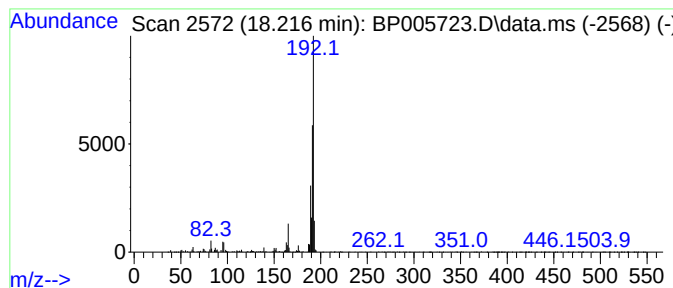
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	5.12 ng	517272	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
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Instrument :
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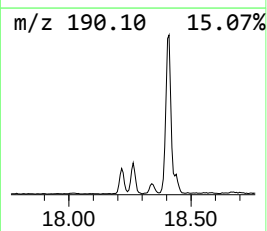
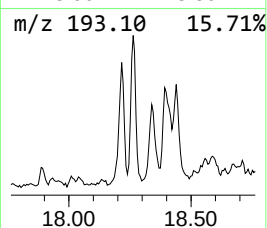
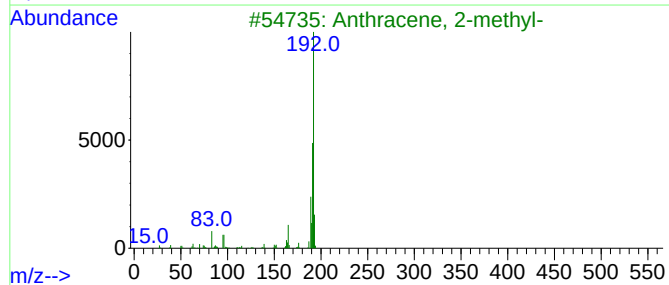
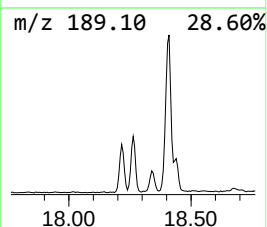
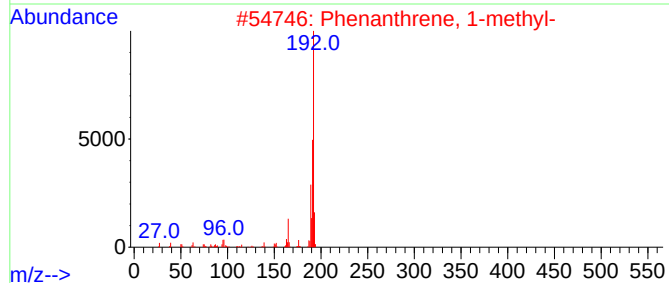
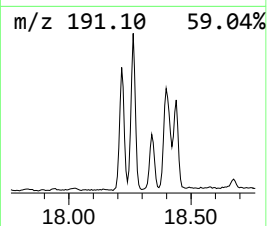
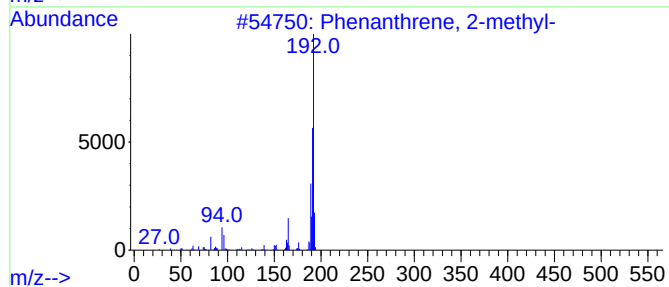
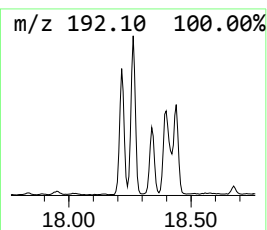
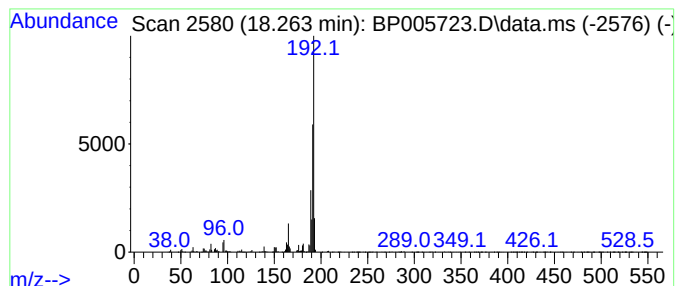
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.263	6.63 ng	670269	Phenanthrene-d10	17.357

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	97
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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18-B10(4-6)

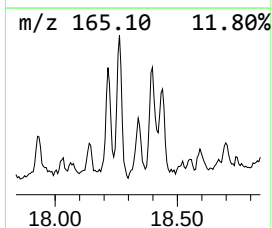
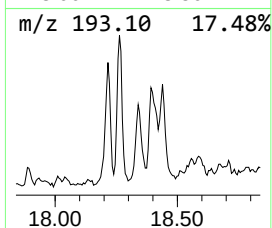
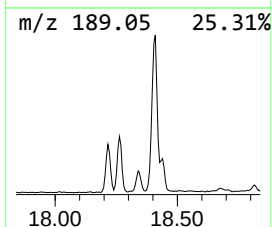
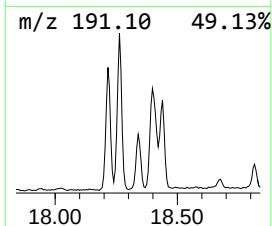
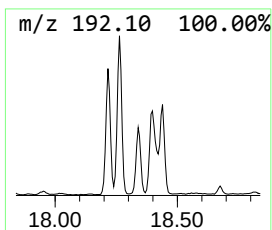
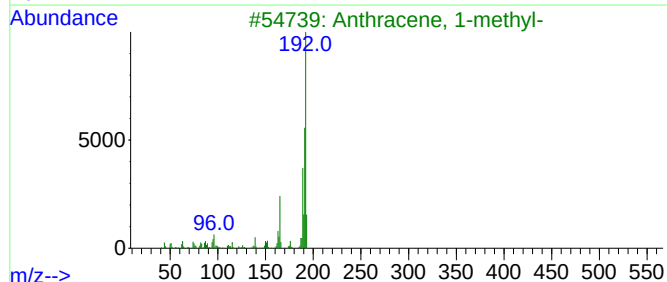
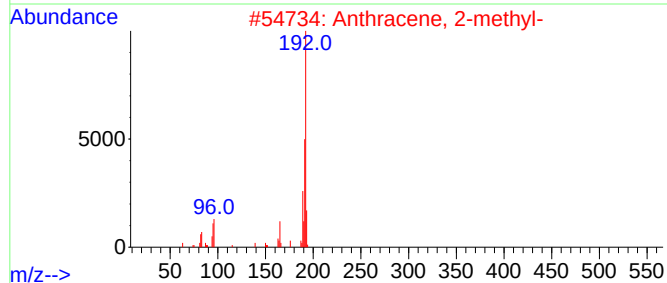
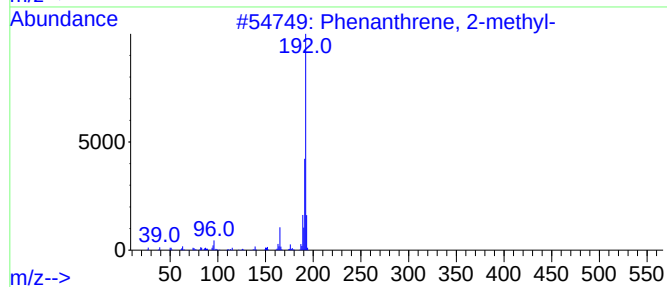
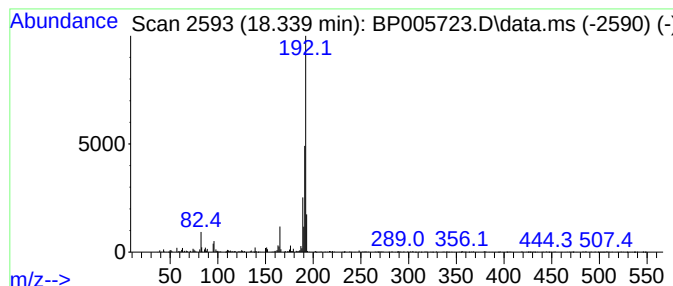
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.339	2.08 ng	209668	Phenanthrene-d10	17.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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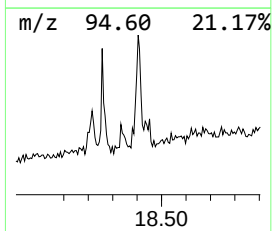
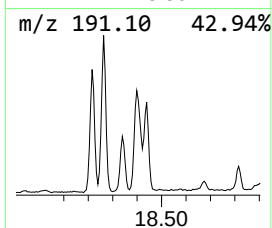
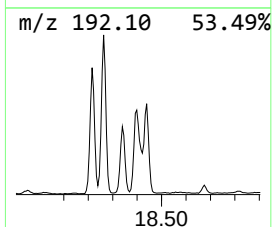
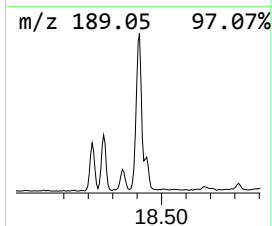
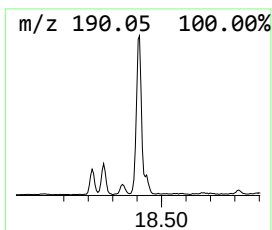
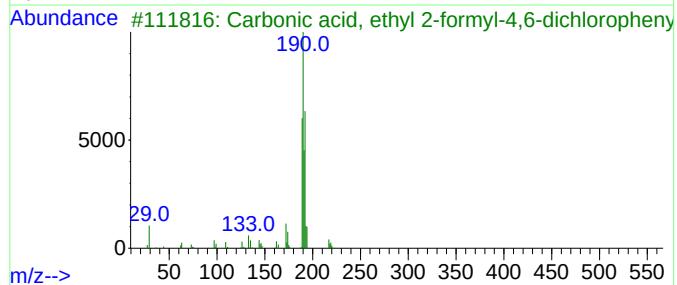
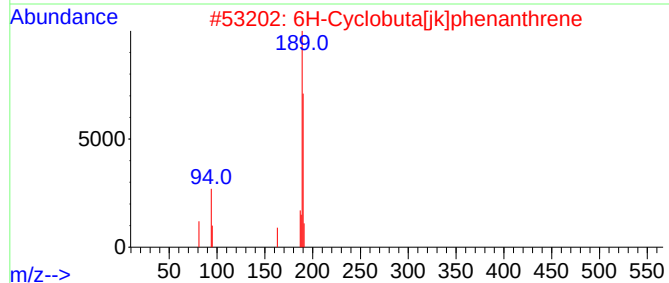
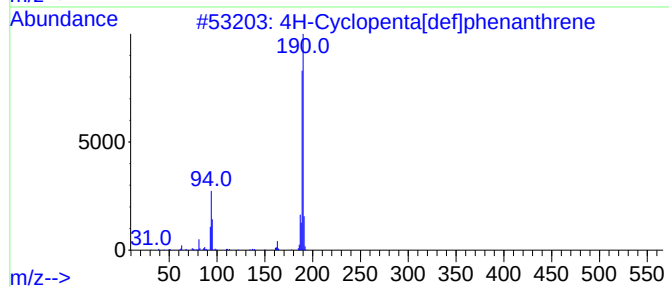
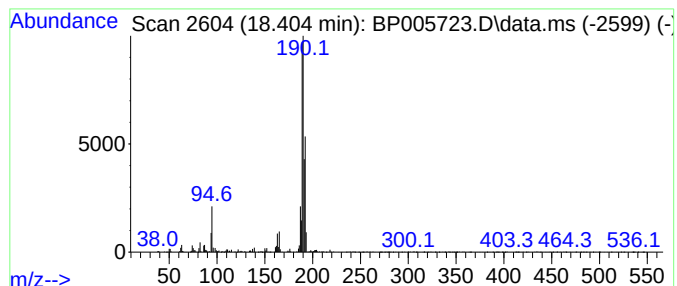
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	9.60 ng	969549	Phenanthrene-d10	17.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
3			Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
4			Methyl diselenide	190	C2H6Se2	007101-31-7	41
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
Sample : M2589-13 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
18-B10(4-6)

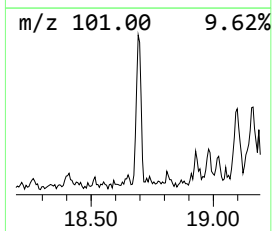
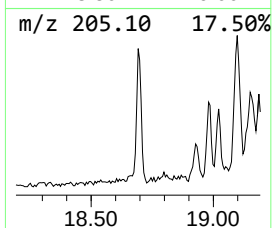
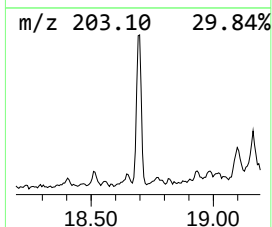
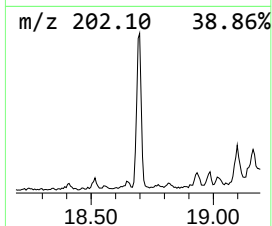
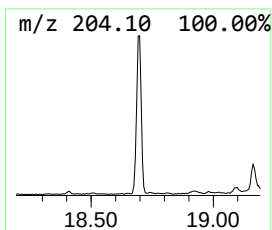
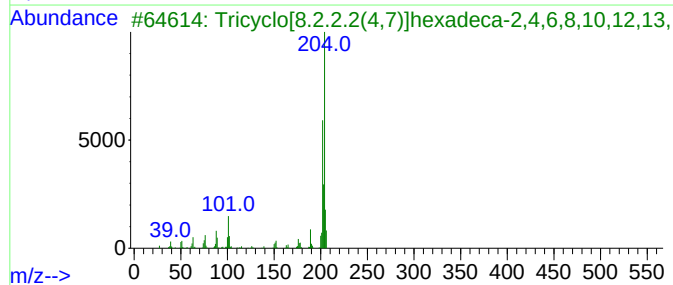
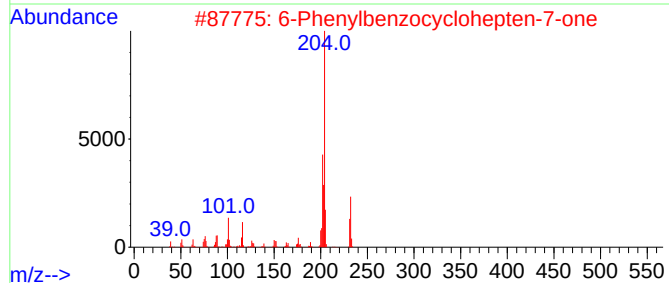
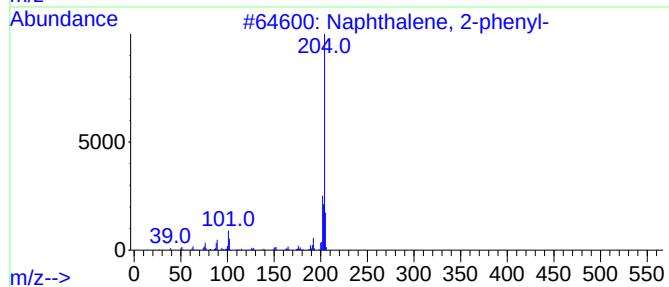
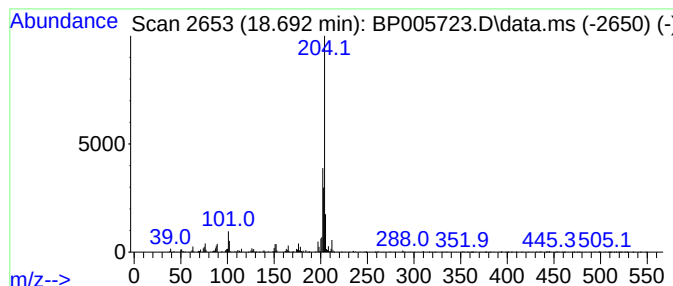
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.692	2.79 ng	281954	Phenanthrene-d10	17.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
2		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	78
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
4		9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	50
5		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
Sample : M2589-13 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

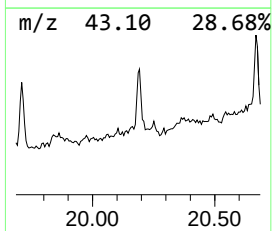
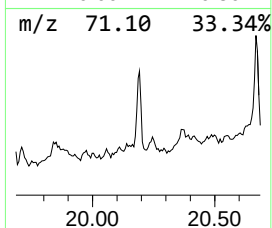
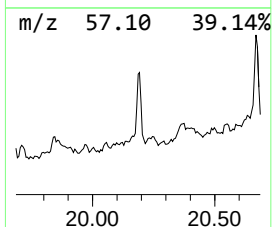
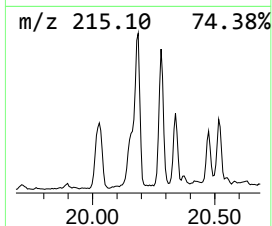
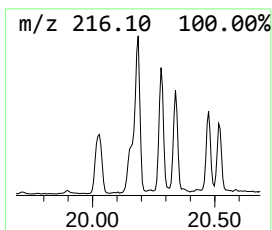
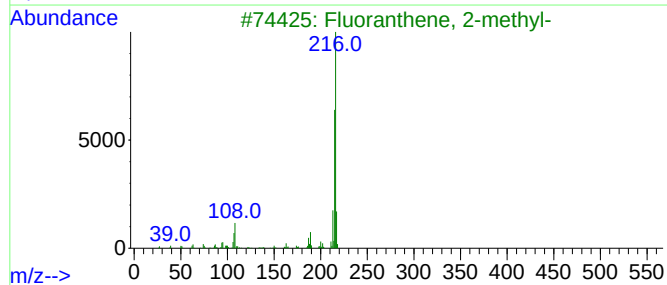
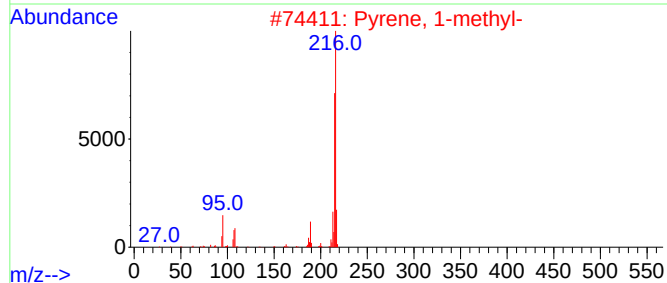
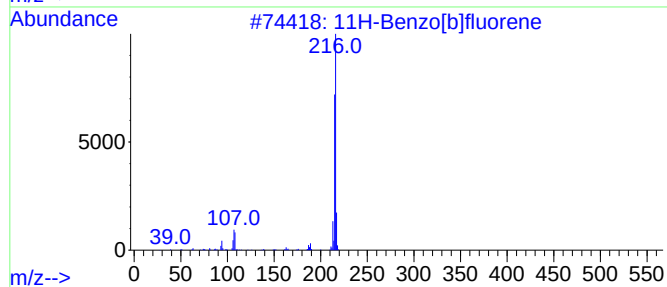
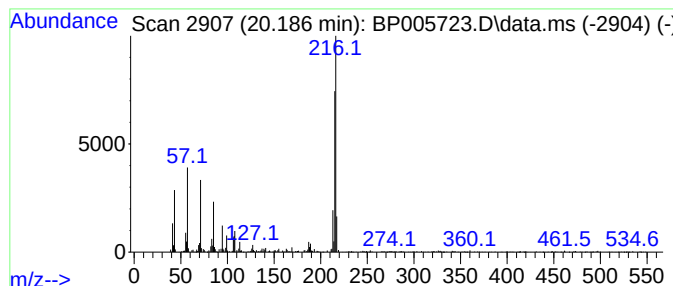
Instrument :
BNA_P
ClientSampleId :
18-B10(4-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.186	3.21 ng	801194	Chrysene-d12	21.422	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060421\
Data File : BP005723.D
Acq On : 04 Jun 2021 21:42
Operator : CG/JU
Sample : M2589-13 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B10(4-6)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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
2-Pentanone, 4-...	5.075	8.3	ng	394155	1	7.987	950816	20.0
(1R)-2,6,6-Trim...	6.663	3.0	ng	142021	1	7.987	950816	20.0
Heptadecane, 2,...	9.216	2.0	ng	95427	1	7.987	950816	20.0
Tetradecane	13.440	2.3	ng	199542	3	14.610	1735360	20.0
Naphthalene, 2,...	13.934	2.3	ng	199078	3	14.610	1735360	20.0
9H-Fluorene, 1-...	16.651	2.1	ng	215179	4	17.357	2020510	20.0
Naphtho[1,2-b]t...	17.169	3.5	ng	356050	4	17.357	2020510	20.0
Phenanthrene, 1...	18.216	5.1	ng	517272	4	17.357	2020510	20.0
Phenanthrene, 2...	18.263	6.6	ng	670269	4	17.357	2020510	20.0
Anthracene, 2-m...	18.339	2.1	ng	209668	4	17.357	2020510	20.0
4H-Cyclopenta[d...	18.404	9.6	ng	969549	4	17.357	2020510	20.0
Naphthalene, 2-...	18.692	2.8	ng	281954	4	17.357	2020510	20.0
11H-Benzo[b]flu...	20.186	3.2	ng	801194	5	21.422	4986660	20.0