

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
 Data File : BM029665.D
 Acq On : 27 Apr 2021 00:27
 Operator : CG/JU
 Sample : M2176-18 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 STOCKPILE-COMP-4

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM029665.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.216	390	396	405	rBV	456173	664247	19.91%	2.013%
2	6.310	577	582	589	rVB2	40616	62806	1.88%	0.190%
3	6.798	657	665	676	rBV	422167	689151	20.65%	2.088%
4	7.616	798	804	814	rBV	288657	449922	13.48%	1.363%
5	8.769	994	1000	1010	rBV	242711	423692	12.70%	1.284%
6	10.398	1269	1277	1283	rBV	359505	609552	18.27%	1.847%
7	10.451	1283	1286	1292	rVB	22924	36888	1.11%	0.112%
8	12.880	1692	1699	1712	rBV	773690	1116503	33.46%	3.384%
9	13.727	1837	1843	1847	rBV2	22784	36194	1.08%	0.110%
10	13.986	1881	1887	1899	rBV	134084	223368	6.69%	0.677%
11	14.110	1902	1908	1915	rVV2	17395	40971	1.23%	0.124%
12	14.263	1927	1934	1941	rBV	568578	874099	26.20%	2.649%
13	14.327	1941	1945	1954	rVB	39739	67549	2.02%	0.205%
14	14.668	1997	2003	2012	rVB	41492	67616	2.03%	0.205%
15	15.315	2108	2113	2118	rBV	89994	127274	3.81%	0.386%
16	15.610	2159	2163	2168	rBV	29396	46539	1.39%	0.141%
17	15.757	2180	2188	2197	rBV	659217	935755	28.04%	2.836%
18	15.998	2222	2229	2236	rVB6	34758	71800	2.15%	0.218%
19	16.321	2277	2284	2289	rBV2	28350	60887	1.82%	0.185%
20	16.668	2339	2343	2349	rVB2	33255	49805	1.49%	0.151%
21	16.745	2352	2356	2360	rBV3	26592	46105	1.38%	0.140%
22	16.833	2367	2371	2379	rVB	55213	94897	2.84%	0.288%
23	17.009	2395	2401	2405	rBV	735104	1063741	31.88%	3.224%
24	17.051	2405	2408	2415	rVV	912037	1229058	36.83%	3.725%
25	17.145	2419	2424	2430	rVV	323437	475839	14.26%	1.442%
26	17.204	2430	2434	2440	rVB2	30712	57109	1.71%	0.173%
27	17.415	2461	2470	2474	rBV	45930	87236	2.61%	0.264%
28	17.527	2485	2489	2495	rVB4	32021	48654	1.46%	0.147%
29	17.592	2496	2500	2507	rVB5	25382	40511	1.21%	0.123%
30	17.886	2545	2550	2553	rBV	125521	177576	5.32%	0.538%
31	17.933	2553	2558	2567	rVV	195897	372822	11.17%	1.130%
32	18.015	2567	2572	2576	rVV	89168	132900	3.98%	0.403%
33	18.080	2577	2583	2587	rVV2	248529	450275	13.49%	1.365%
34	18.115	2587	2589	2596	rVB	90882	118032	3.54%	0.358%
35	18.392	2632	2636	2639	rBV	90082	113598	3.40%	0.344%
36	18.433	2639	2643	2649	rVB3	49690	69507	2.08%	0.211%

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Filtering: 5

Sampling : 1

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Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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37	18.639	2675	2678	2682	rVB2	29366	33701	1.01%	0.102%
38	18.686	2683	2686	2689	rVV	39878	48915	1.47%	0.148%
39	18.727	2690	2693	2697	rVB2	36934	45394	1.36%	0.138%
40	18.809	2702	2707	2711	rBV	103794	181518	5.44%	0.550%
41	18.950	2727	2731	2739	rVB2	101091	173342	5.19%	0.525%
42	19.074	2746	2752	2760	rBV	2240923	2852090	85.47%	8.643%
43	19.174	2766	2769	2771	rBV2	35583	39264	1.18%	0.119%
44	19.221	2774	2777	2781	rVB	98888	141558	4.24%	0.429%
45	19.433	2809	2813	2822	rVB	1777674	2529158	75.80%	7.664%
46	19.509	2823	2826	2834	rVB2	85637	136274	4.08%	0.413%
47	19.645	2841	2849	2853	rBV	1359720	1774940	53.19%	5.379%
48	19.762	2864	2869	2876	rBV2	120317	300720	9.01%	0.911%
49	19.903	2888	2893	2894	rBV4	91381	150400	4.51%	0.456%
50	19.933	2894	2898	2903	rVB2	331867	448199	13.43%	1.358%
51	20.033	2911	2915	2919	rBV	260102	321727	9.64%	0.975%
52	20.092	2920	2925	2929	rVV2	179984	275489	8.26%	0.835%
53	20.227	2946	2948	2952	rVB	88742	93507	2.80%	0.283%
54	20.274	2953	2956	2961	rVV2	83787	120135	3.60%	0.364%
55	20.680	3022	3025	3031	rVB	148342	214072	6.42%	0.649%
56	20.844	3050	3053	3056	rVB	114750	139595	4.18%	0.423%
57	20.880	3057	3059	3062	rVB	136230	131058	3.93%	0.397%
58	20.921	3062	3066	3070	rBV2	299194	408822	12.25%	1.239%
59	21.186	3106	3111	3116	rBV2	1681207	3336815	100.00%	10.112%
60	21.233	3116	3119	3127	rVB	1076623	1392528	41.73%	4.220%
61	21.350	3136	3139	3144	rBV2	181495	301490	9.04%	0.914%
62	22.733	3369	3374	3379	rBV	1048286	2216048	66.41%	6.716%
63	23.197	3449	3453	3462	rVB	581781	1069988	32.07%	3.243%
64	23.291	3464	3469	3477	rVB	983448	1723732	51.66%	5.224%
65	23.385	3480	3485	3490	rBV2	699862	1165411	34.93%	3.532%

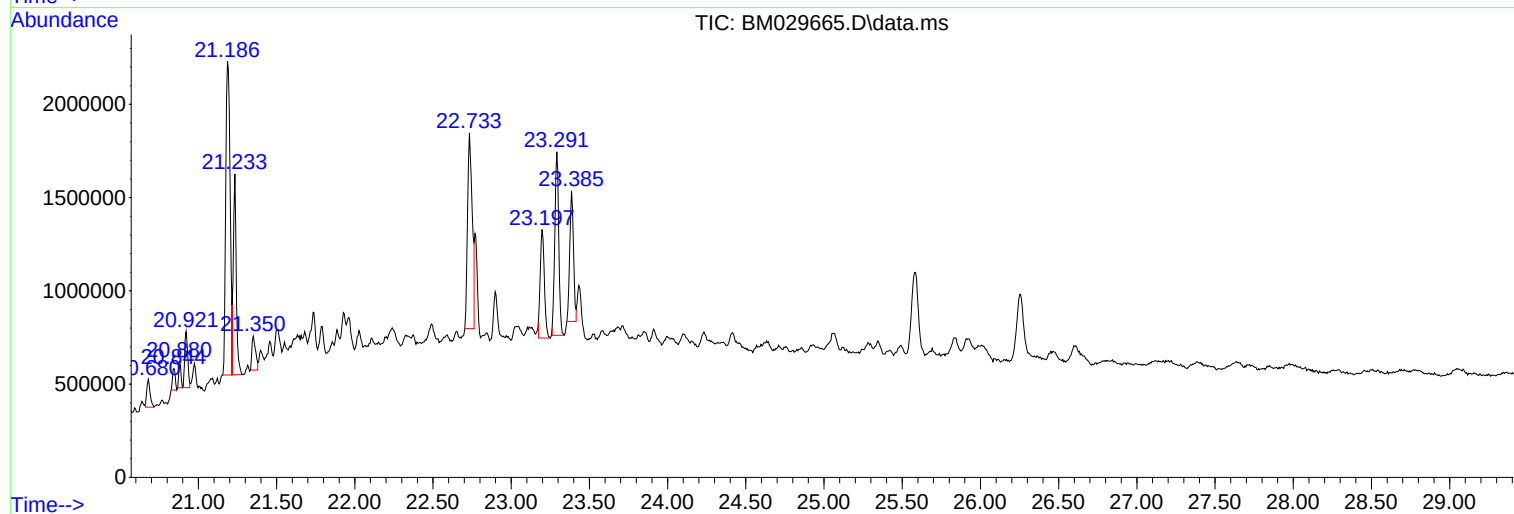
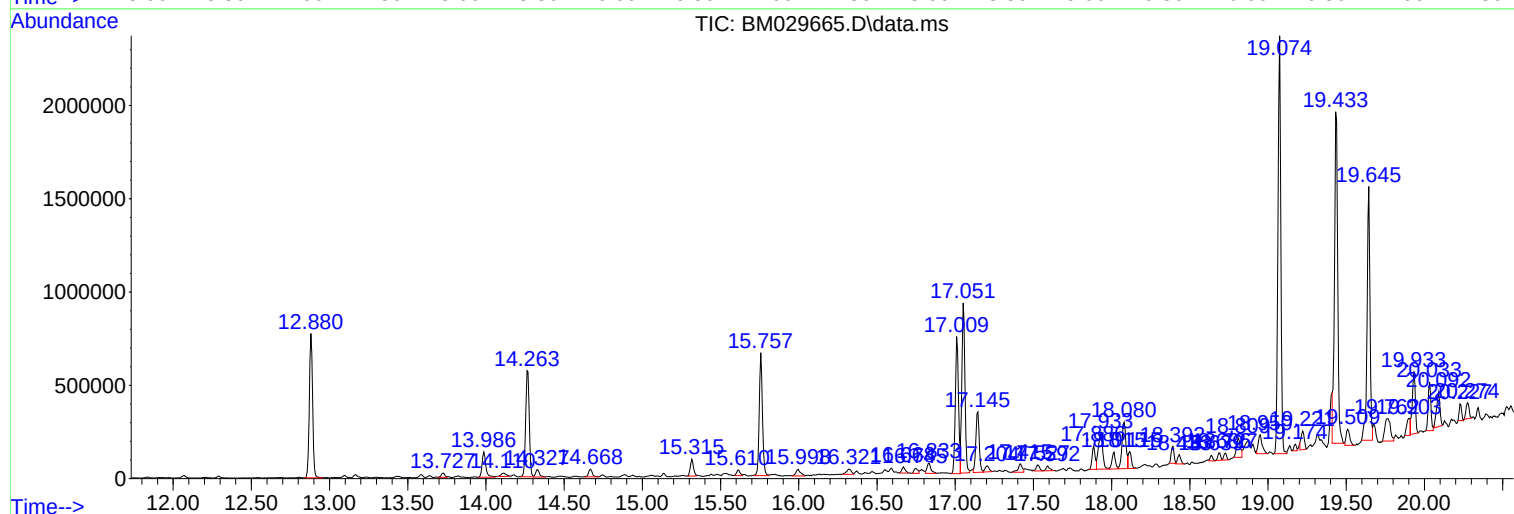
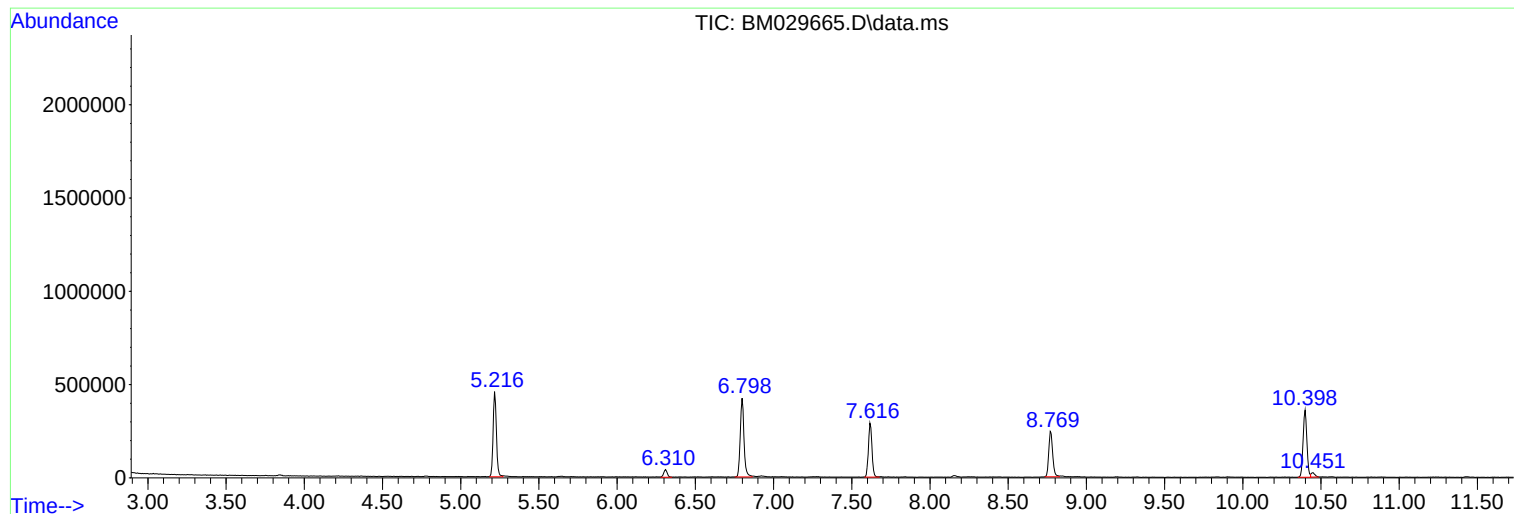
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
STOCKPILE-COMP-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.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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Data File : BM029665.D
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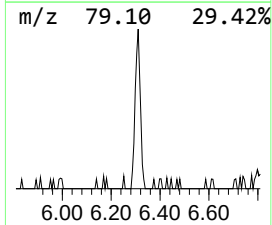
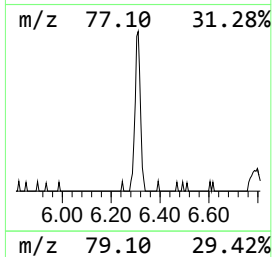
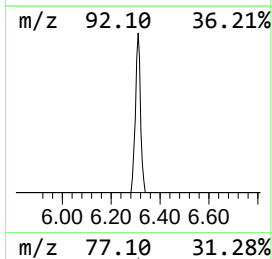
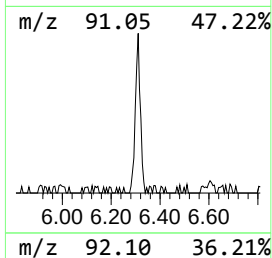
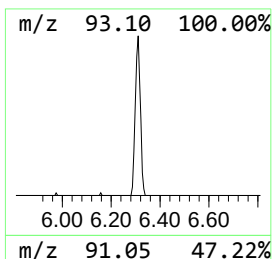
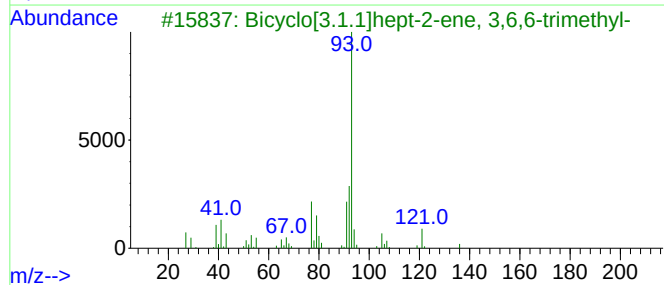
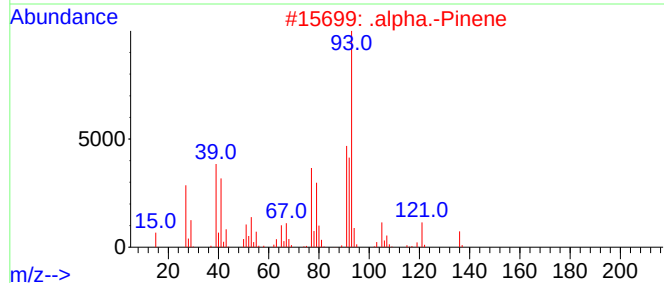
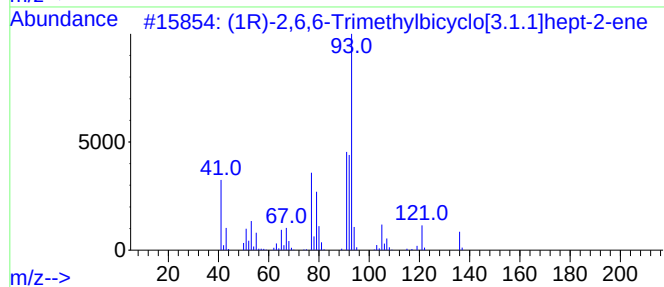
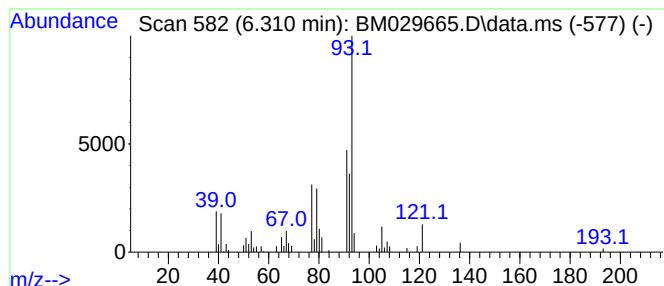
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.310	2.79 ng	62806	1,4-Dichlorobenzene-d4	7.616

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	95
3		Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91
4		Tricyclo[2.2.1.0(2,6)]heptane, 1,1,1-trimethyl-	136	C10H16	000508-32-7	91
5		(+)-3-Carene	136	C10H16	000498-15-7	91



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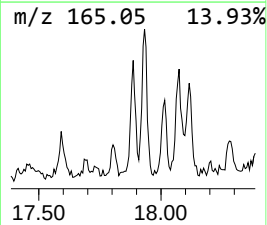
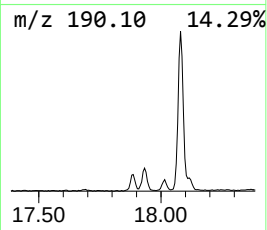
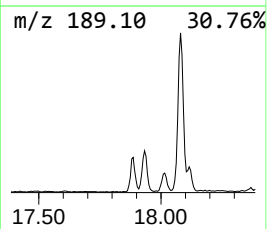
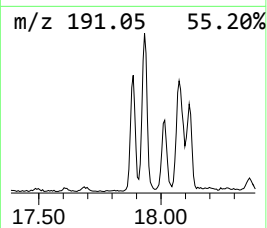
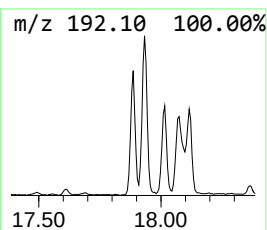
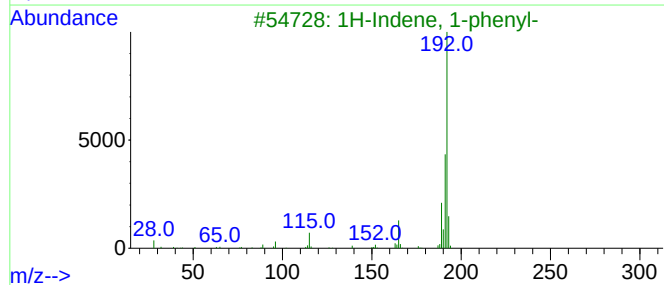
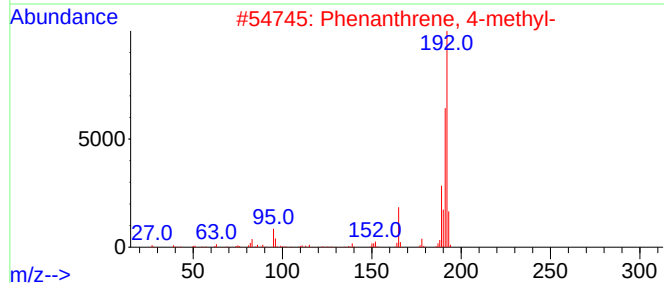
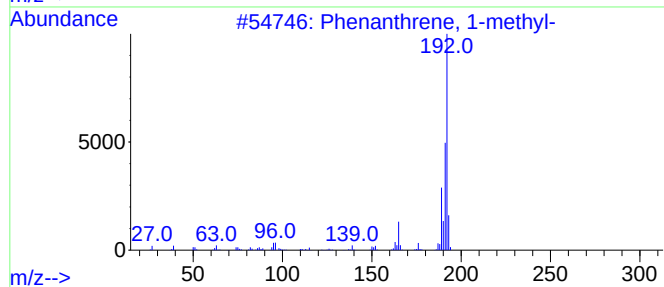
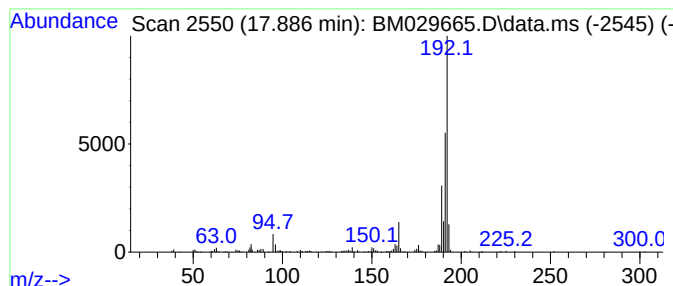
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.886	3.34 ng	177576	Phenanthrene-d10	17.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
3		1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	94
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



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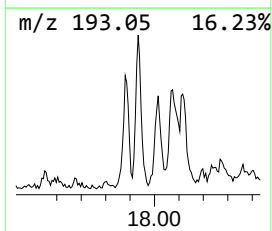
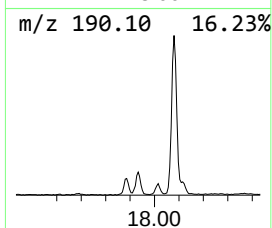
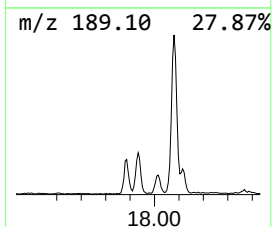
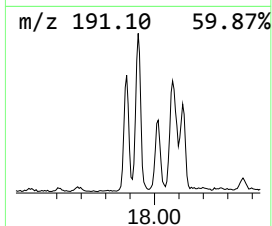
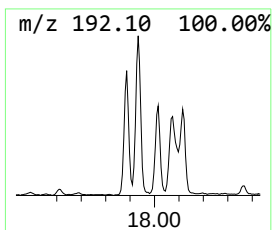
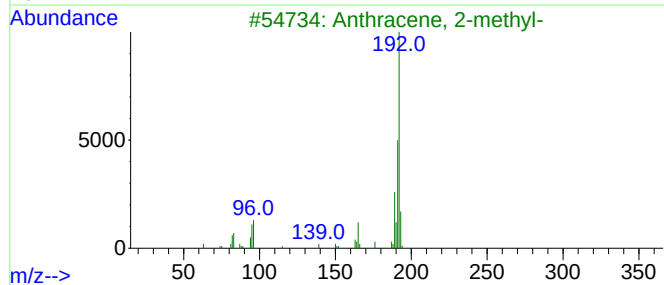
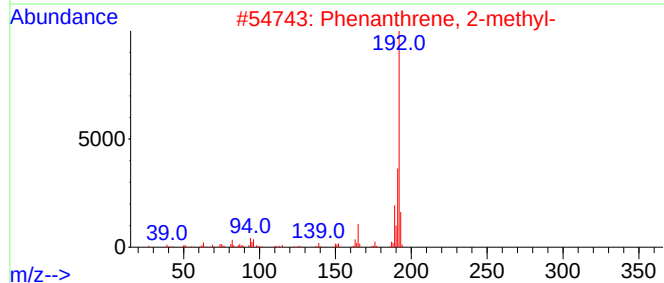
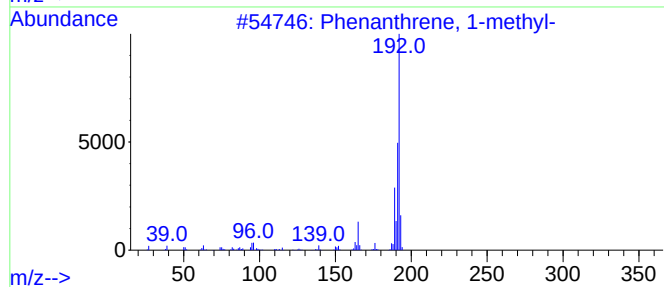
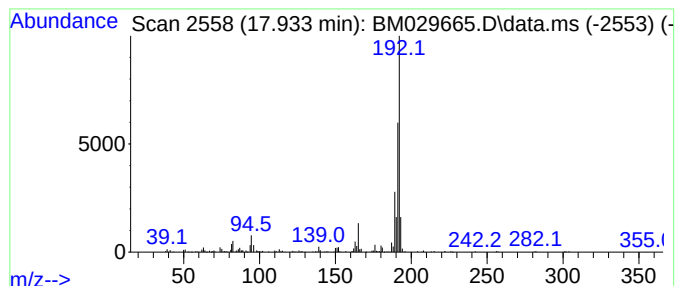
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.933	7.01 ng	372822	Phenanthrene-d10	17.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



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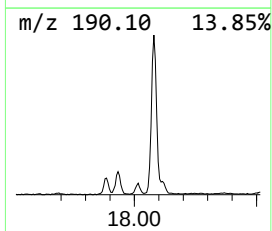
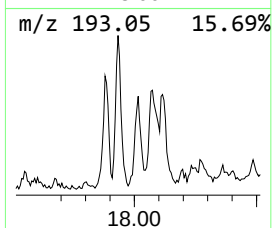
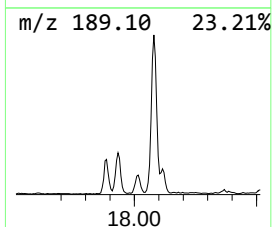
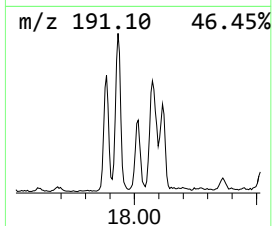
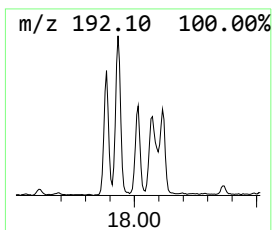
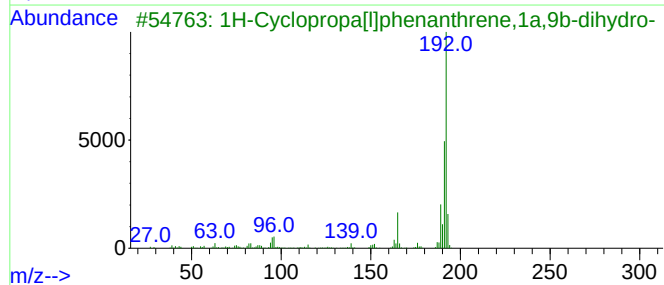
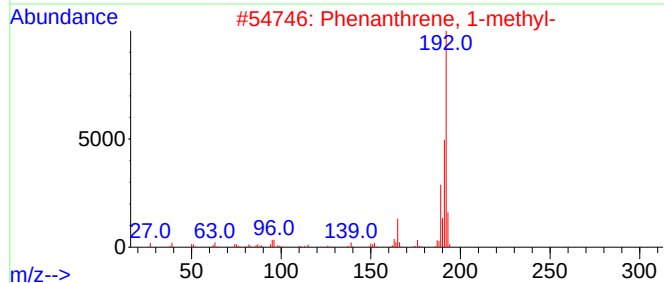
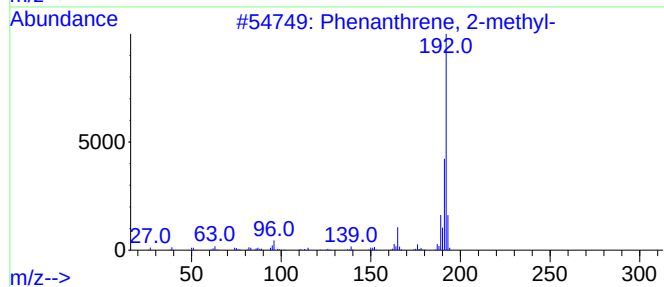
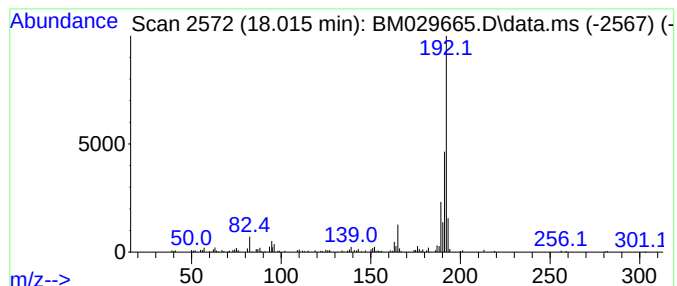
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.015	2.50 ng	132900	Phenanthrene-d10	17.009

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			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
STOCKPILE-COMP-4

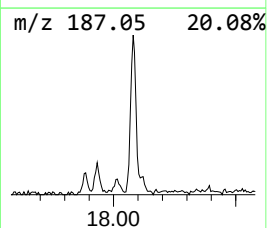
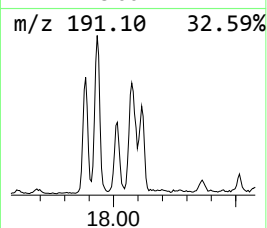
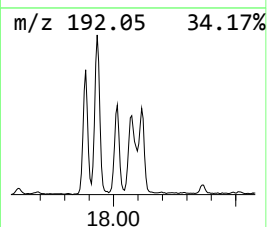
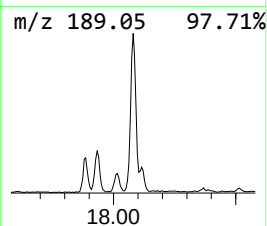
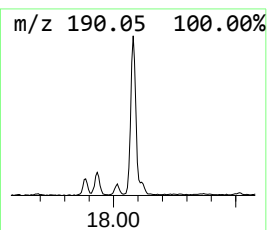
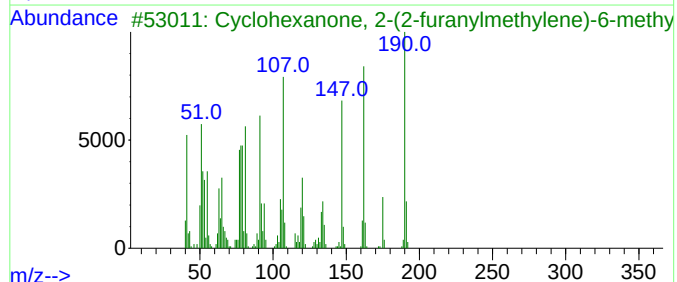
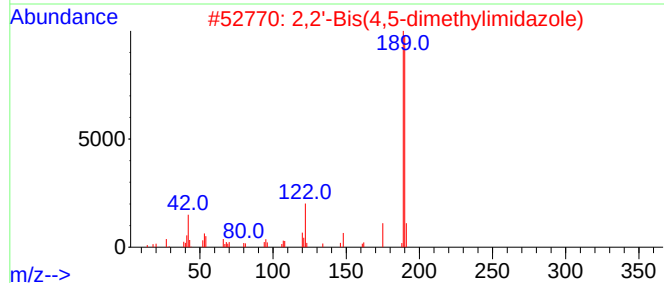
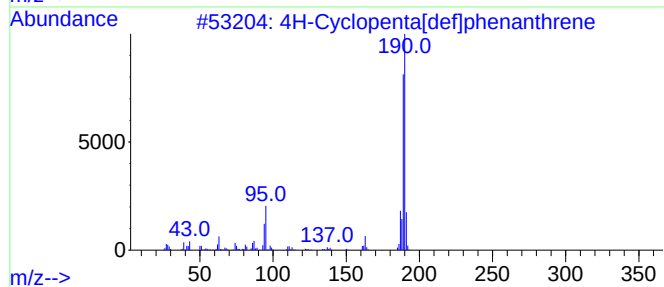
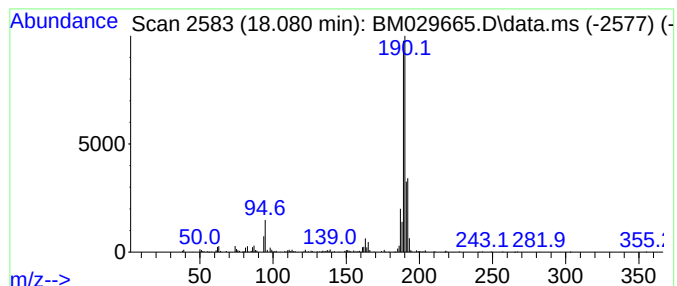
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.080	8.47 ng	450275	Phenanthrene-d10	17.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
3			Cyclohexanone, 2-(2-furanylmethy...	190	C12H14O2	056053-07-7	27
4			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5			4-Isothiazolecarbonitrile, 5-chl...	190	C5H3ClN2S2	054798-93-5	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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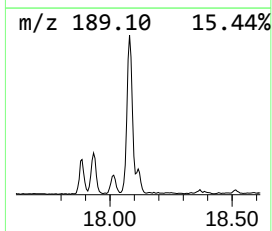
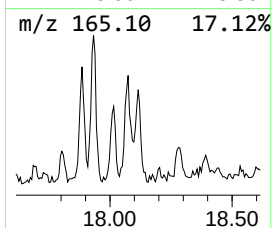
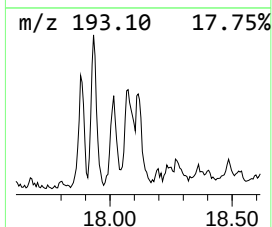
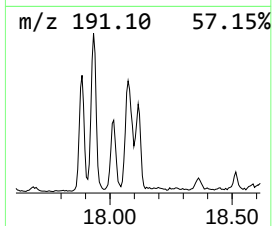
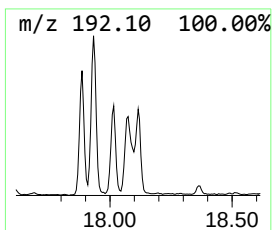
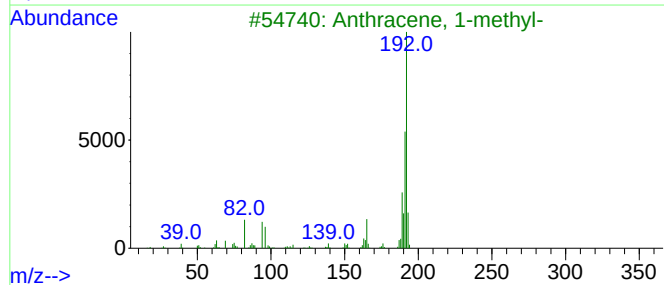
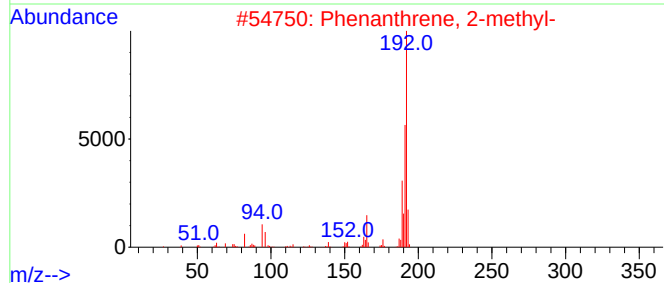
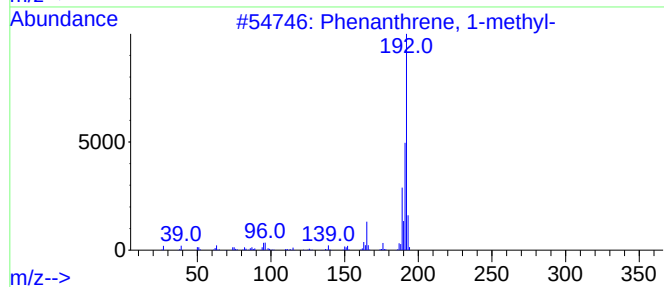
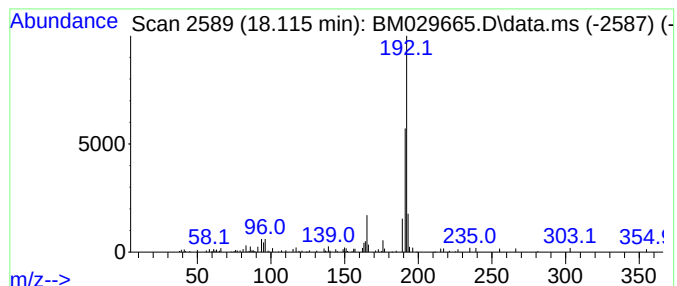
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.115	2.22 ng	118032	Phenanthrene-d10	17.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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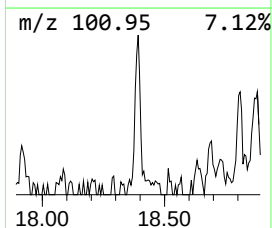
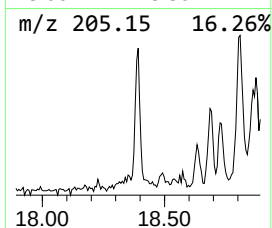
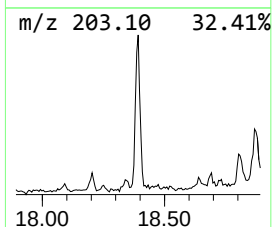
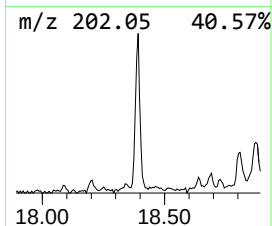
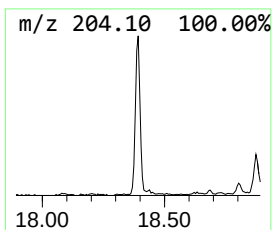
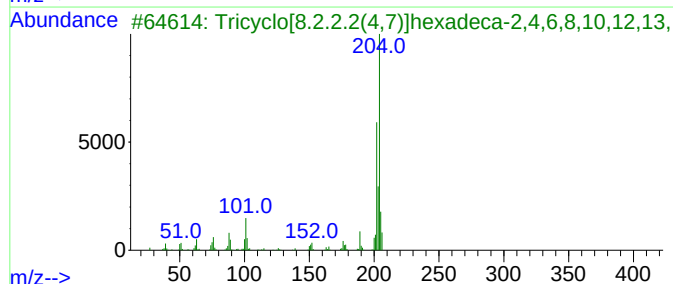
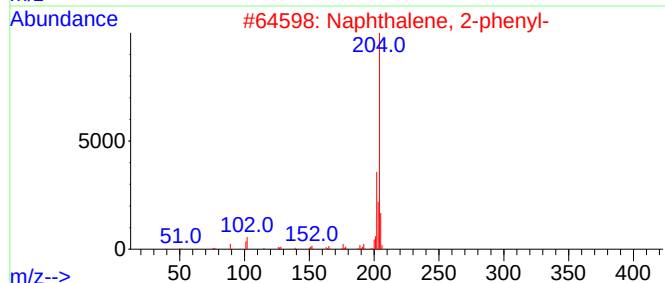
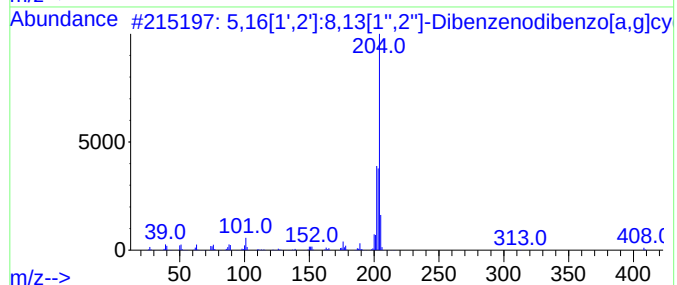
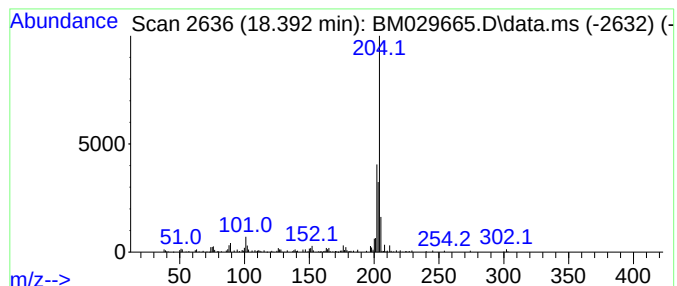
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.392	2.14 ng	113598	Phenanthrene-d10	17.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:	8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86	
2	Naphthalene, 2-phenyl-		204	C16H12	000612-94-2	70	
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...		204	C16H12	006572-60-7	62	
4	3,4-Biphenyldicarbonitrile		204	C14H8N2	004128-63-6	47	
5	9-Acridinecarbonitrile		204	C14H8N2	005326-19-2	47	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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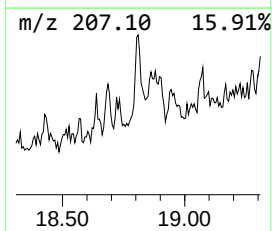
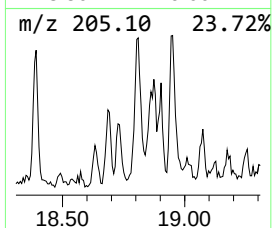
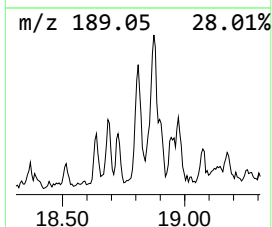
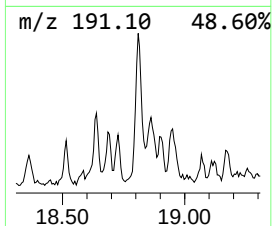
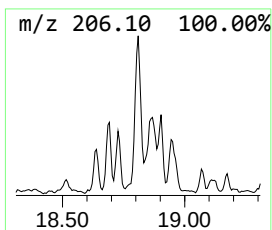
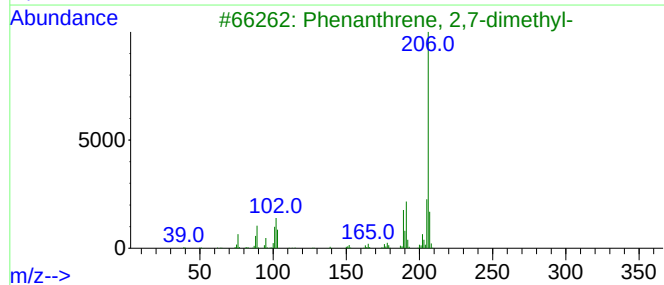
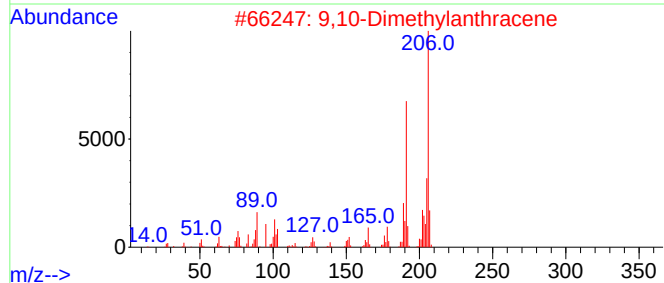
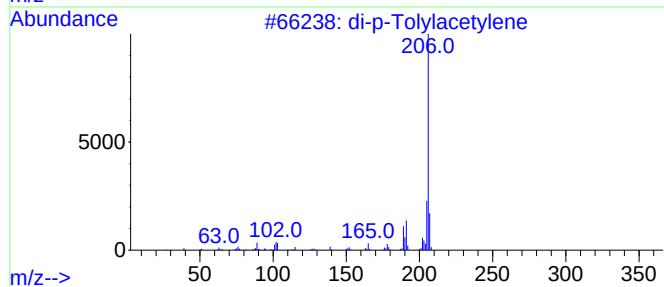
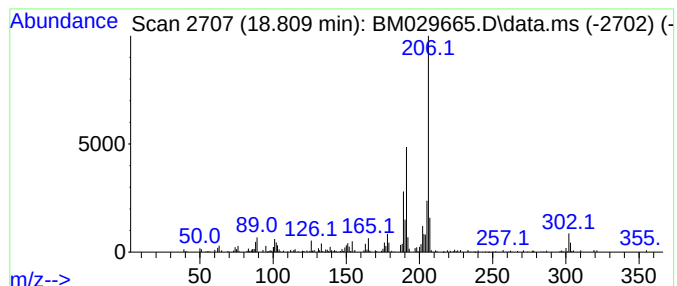
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 di-p-Tolylacetylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.809	3.41 ng	181518	Phenanthrene-d10	17.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-p-Tolylacetylene	206	C16H14	002789-88-0	91
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	87
4			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	83
5			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
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ALS Vial : 18 Sample Multiplier: 1

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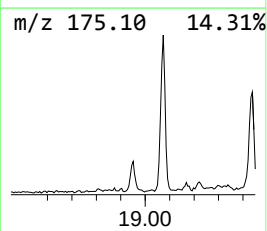
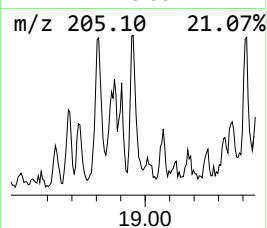
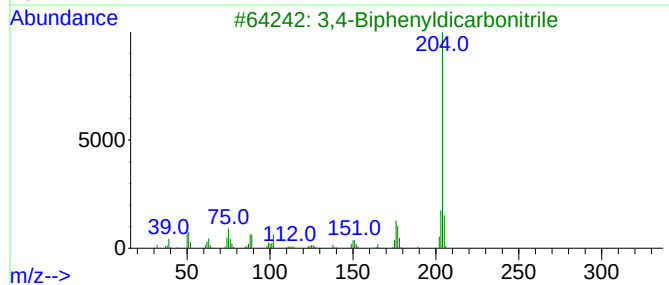
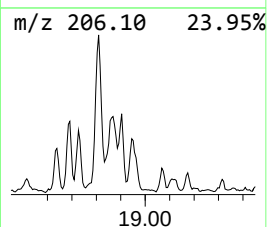
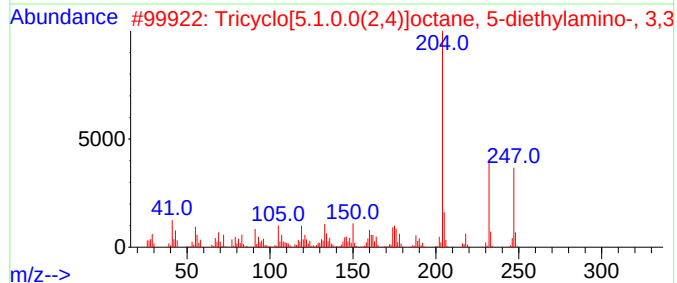
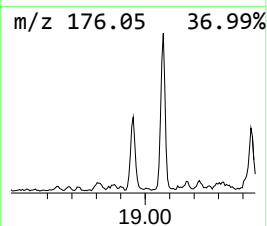
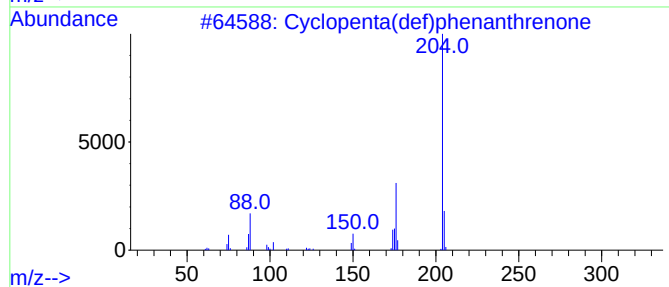
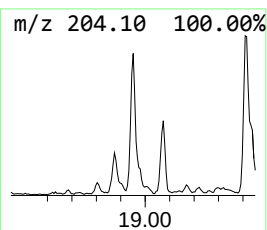
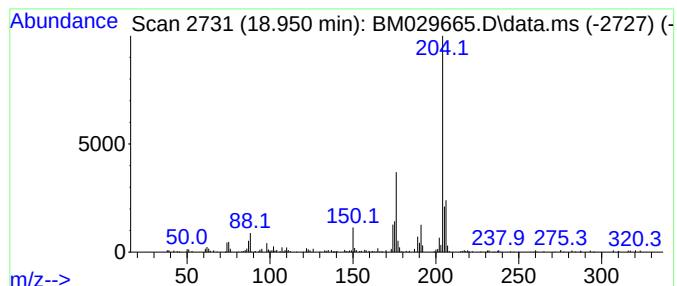
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.951	3.26 ng	173342	Phenanthrene-d10	17.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	81
2			Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	59
3			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	59
4			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	47
5			[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
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Acq On : 27 Apr 2021 00:27
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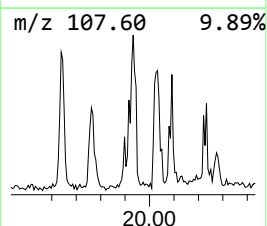
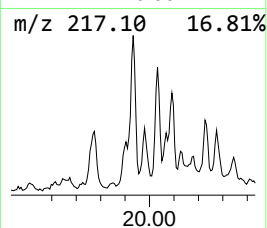
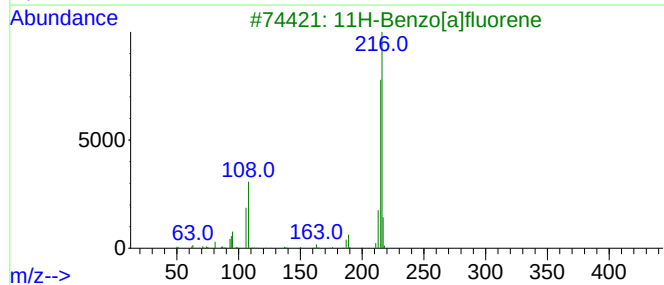
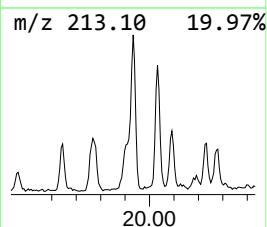
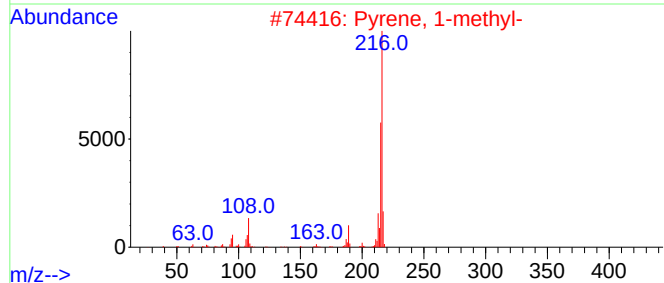
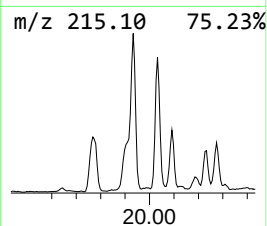
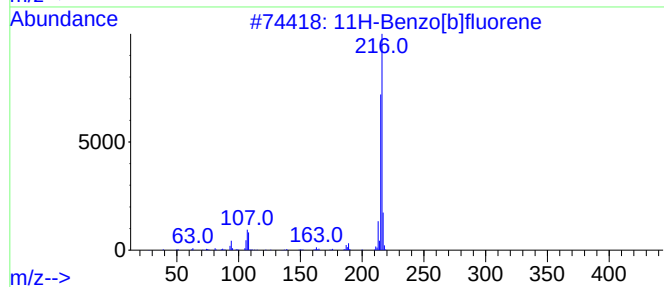
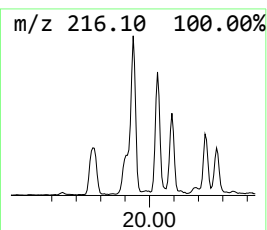
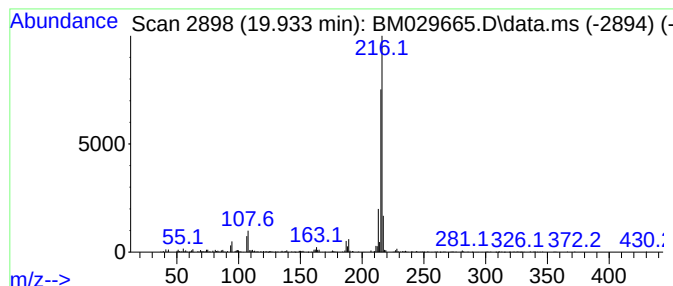
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.933	2.69 ng	448199	Chrysene-d12	21.197

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
5		.beta.-Carboline, 1,2,3,4-tetra...	260	C14H16N2O3	1000124-14-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
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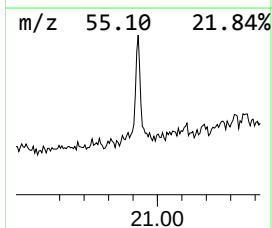
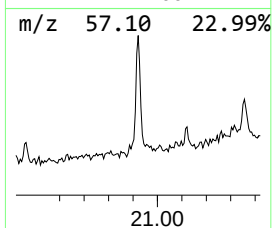
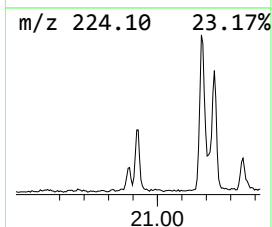
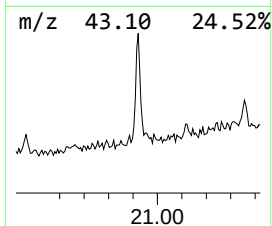
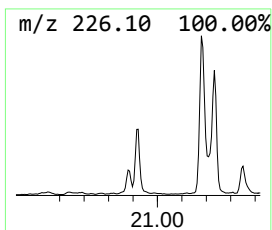
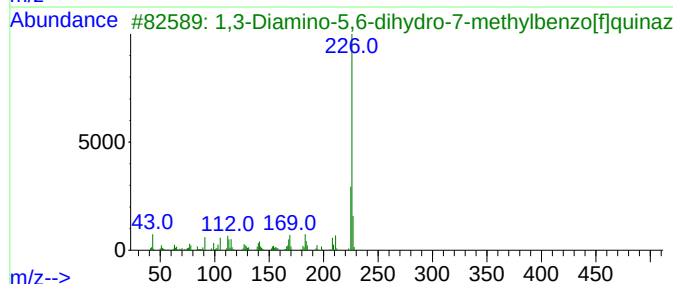
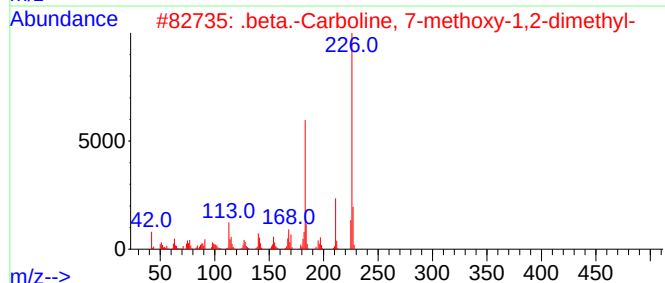
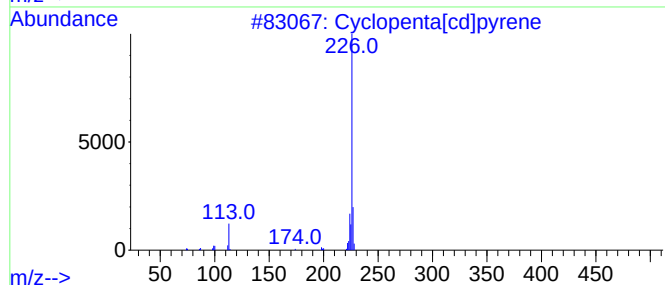
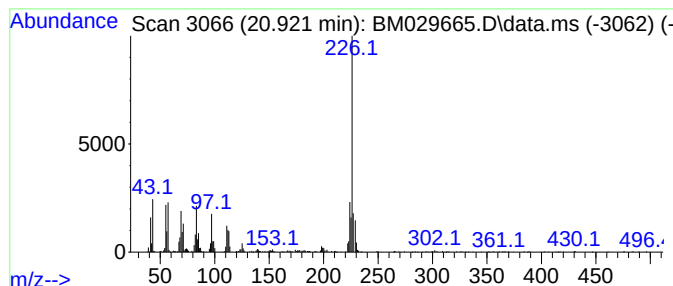
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclopenta[cd]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.921	2.45 ng	408822	Chrysene-d12	21.197

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	53
2			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	50
3			1,3-Diamino-5,6-dihydro-7-methyl...	226	C13H14N4	037436-37-6	50
4			9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	46
5			Anthracene, 1,8-diethynyl-	226	C18H10	078053-58-4	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
STOCKPILE-COMP-4

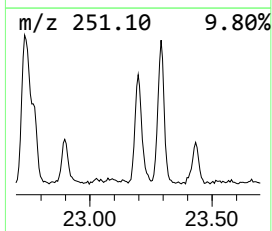
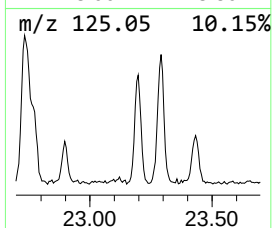
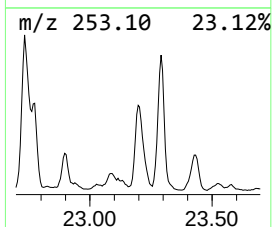
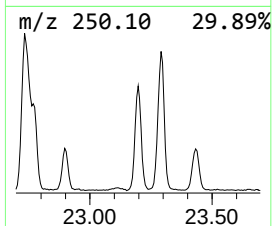
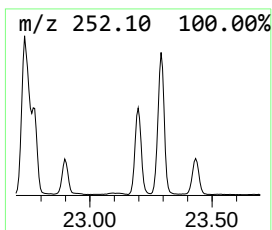
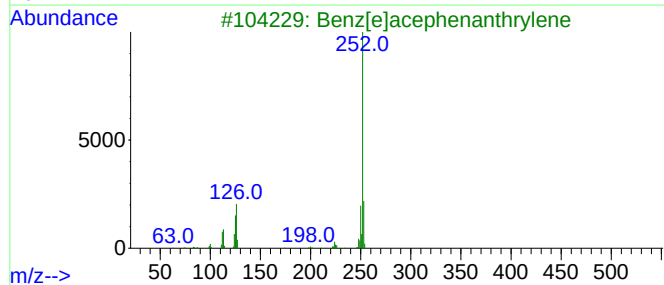
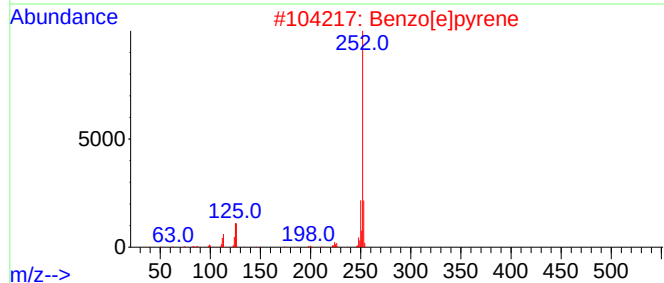
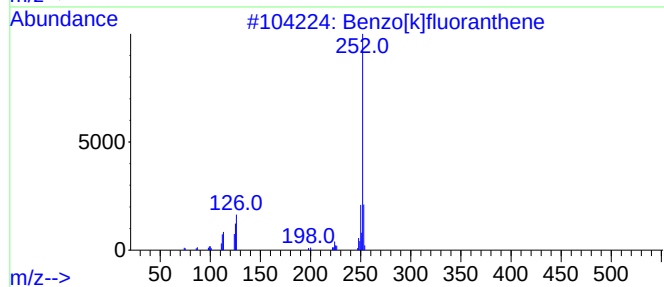
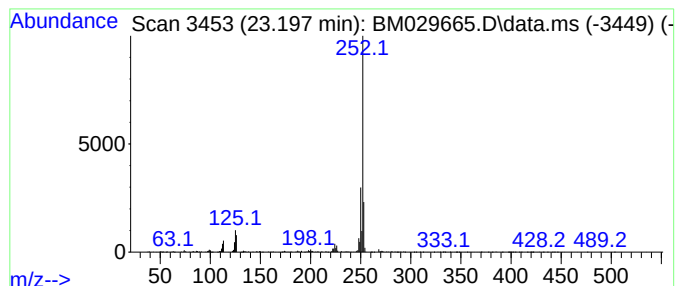
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.197	18.36 ng	1069990	Perylene-d12	23.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
2			Benzo[e]pyrene	252	C20H12	000192-97-2	96
3			Benzo[a]acephenanthrylene	252	C20H12	000205-99-2	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042621\
Data File : BM029665.D
Acq On : 27 Apr 2021 00:27
Operator : CG/JU
Sample : M2176-18 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
STOCKPILE-COMP-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.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
(1R)-2,6,6-Trim...	6.310	2.8	ng	62806	1	7.616	449922	20.0
Phenanthrene, 1...	17.886	3.3	ng	177576	4	17.009	1063740	20.0
Phenanthrene, 2...	17.933	7.0	ng	372822	4	17.009	1063740	20.0
1H-Cyclopropa[l...	18.015	2.5	ng	132900	4	17.009	1063740	20.0
4H-Cyclopenta[d...	18.080	8.5	ng	450275	4	17.009	1063740	20.0
Anthracene, 1-m...	18.115	2.2	ng	118032	4	17.009	1063740	20.0
Naphthalene, 2-...	18.392	2.1	ng	113598	4	17.009	1063740	20.0
di-p-Tolylacety...	18.809	3.4	ng	181518	4	17.009	1063740	20.0
Cyclopenta(def)...	18.951	3.3	ng	173342	4	17.009	1063740	20.0
11H-Benzo[b]flu...	19.933	2.7	ng	448199	5	21.197	3336820	20.0
Cyclopenta[cd]p...	20.921	2.5	ng	408822	5	21.197	3336820	20.0
Benzo[e]pyrene	23.197	18.4	ng	1069990	6	23.386	1165410	20.0