

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
 Data File : BP006106.D
 Acq On : 29 Jun 2021 23:44
 Operator : CG/JU
 Sample : M2891-01 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TR-04-062821

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-BP060821.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.016	323	328	335	rVB2	14549	24612	1.14%	0.079%
2	5.481	399	407	427	rBV	732871	1162219	53.75%	3.746%
3	7.087	668	680	696	rBV	697479	1249450	57.79%	4.027%
4	7.904	808	819	834	rBV	572078	944571	43.69%	3.045%
5	8.446	904	911	917	rBV3	11317	24748	1.14%	0.080%
6	9.069	1010	1017	1030	rBV	432903	766733	35.46%	2.471%
7	9.234	1041	1045	1057	rVB	10136	30576	1.41%	0.099%
8	10.504	1253	1261	1269	rBV4	19131	52880	2.45%	0.170%
9	10.704	1283	1295	1301	rBV	703251	1228290	56.81%	3.959%
10	10.751	1301	1303	1310	rVB	44752	68946	3.19%	0.222%
11	11.728	1463	1469	1472	rBV	16882	30090	1.39%	0.097%
12	12.363	1573	1577	1584	rVB	36168	63366	2.93%	0.204%
13	12.581	1609	1614	1623	rVB2	55105	102179	4.73%	0.329%
14	13.069	1693	1697	1703	rVB4	17971	30187	1.40%	0.097%
15	13.157	1703	1712	1723	rBV	1272129	1813906	83.90%	5.847%
16	13.698	1799	1804	1805	rBV	36019	42707	1.98%	0.138%
17	13.857	1826	1831	1836	rBV	72817	126298	5.84%	0.407%
18	13.910	1836	1840	1844	rVV	33410	46582	2.15%	0.150%
19	13.992	1851	1854	1856	rBV	23211	30815	1.43%	0.099%
20	14.092	1867	1871	1874	rBV	34113	46561	2.15%	0.150%
21	14.257	1894	1899	1908	rBV	137758	212336	9.82%	0.684%
22	14.357	1912	1916	1921	rBV3	19660	25843	1.20%	0.083%
23	14.422	1921	1927	1935	rBV4	23558	55756	2.58%	0.180%
24	14.534	1936	1946	1952	rVV	981062	1469727	67.98%	4.738%
25	14.598	1952	1957	1964	rVB	146669	243363	11.26%	0.784%
26	14.745	1976	1982	1990	rBV5	25898	68091	3.15%	0.219%
27	14.934	2008	2014	2021	rVB	136687	204256	9.45%	0.658%
28	15.010	2022	2027	2031	rVB	38682	52937	2.45%	0.171%
29	15.145	2045	2050	2056	rBV3	43731	88351	4.09%	0.285%
30	15.204	2057	2060	2064	rVB	26354	34041	1.57%	0.110%
31	15.322	2072	2080	2083	rBV6	33177	78403	3.63%	0.253%
32	15.351	2083	2085	2091	rVV4	23408	44993	2.08%	0.145%
33	15.404	2091	2094	2100	rVB	24531	35648	1.65%	0.115%
34	15.581	2118	2124	2136	rVB	248041	408643	18.90%	1.317%
35	15.739	2148	2151	2155	rVB4	31888	42328	1.96%	0.136%
36	15.786	2155	2159	2163	rBV4	34533	53641	2.48%	0.173%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
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37	15.875	2170	2174	2180	rVB2	43570	78851	3.65%	0.254%
38	16.028	2191	2200	2208	rBV	836363	1326421	61.35%	4.276%
39	16.257	2233	2239	2248	rVB3	65958	133008	6.15%	0.429%
40	16.580	2288	2294	2299	rBV2	57987	120088	5.55%	0.387%
41	16.633	2299	2303	2309	rVB2	49072	67164	3.11%	0.216%
42	16.733	2316	2320	2325	rVB3	38249	60102	2.78%	0.194%
43	16.939	2352	2355	2360	rVB4	21323	32772	1.52%	0.106%
44	17.098	2378	2382	2390	rVB	78260	143039	6.62%	0.461%
45	17.280	2407	2413	2417	rBV	998032	1435199	66.38%	4.626%
46	17.322	2417	2420	2427	rVV	860801	1122136	51.90%	3.617%
47	17.416	2431	2436	2441	rVV	249667	344488	15.93%	1.110%
48	17.480	2442	2447	2451	rVB3	30940	62103	2.87%	0.200%
49	17.692	2479	2483	2488	rBV2	61503	101010	4.67%	0.326%
50	17.863	2508	2512	2519	rVB2	40641	71212	3.29%	0.230%
51	17.939	2521	2525	2531	rVB	482182	583369	26.98%	1.880%
52	18.145	2556	2560	2565	rBV	197564	365998	16.93%	1.180%
53	18.192	2565	2568	2575	rVB	224063	300224	13.89%	0.968%
54	18.275	2578	2582	2587	rBV2	82013	126713	5.86%	0.408%
55	18.333	2587	2592	2596	rVV2	251232	463980	21.46%	1.496%
56	18.369	2596	2598	2606	rVB	153927	170293	7.88%	0.549%
57	18.439	2608	2610	2614	rBV3	29120	34519	1.60%	0.111%
58	18.627	2639	2642	2646	rVB	82404	101214	4.68%	0.326%
59	18.916	2688	2691	2694	rBV	68811	74266	3.43%	0.239%
60	19.039	2706	2712	2715	rBV2	1687292	2162072	100.00%	6.969%
61	19.069	2715	2717	2729	rVB4	1520432	2019987	93.43%	6.511%
62	19.198	2736	2739	2746	rVB	392718	458259	21.20%	1.477%
63	19.286	2749	2754	2759	rBV	826109	1082959	50.09%	3.491%
64	19.433	2776	2779	2783	rVB	73072	85923	3.97%	0.277%
65	19.610	2805	2809	2810	rBV2	144160	181677	8.40%	0.586%
66	19.639	2810	2814	2822	rVV	871392	1319444	61.03%	4.253%
67	19.710	2823	2826	2832	rVB	115640	191711	8.87%	0.618%
68	19.827	2842	2846	2850	rVB	1535640	1842657	85.23%	5.940%
69	20.216	2909	2912	2916	rBV	161481	186227	8.61%	0.600%
70	20.416	2942	2946	2949	rBV3	218675	340991	15.77%	1.099%
71	21.357	3100	3106	3110	rBV2	1156430	1794469	83.00%	5.784%
72	23.762	3512	3515	3521	rVB2	620819	1034609	47.85%	3.335%

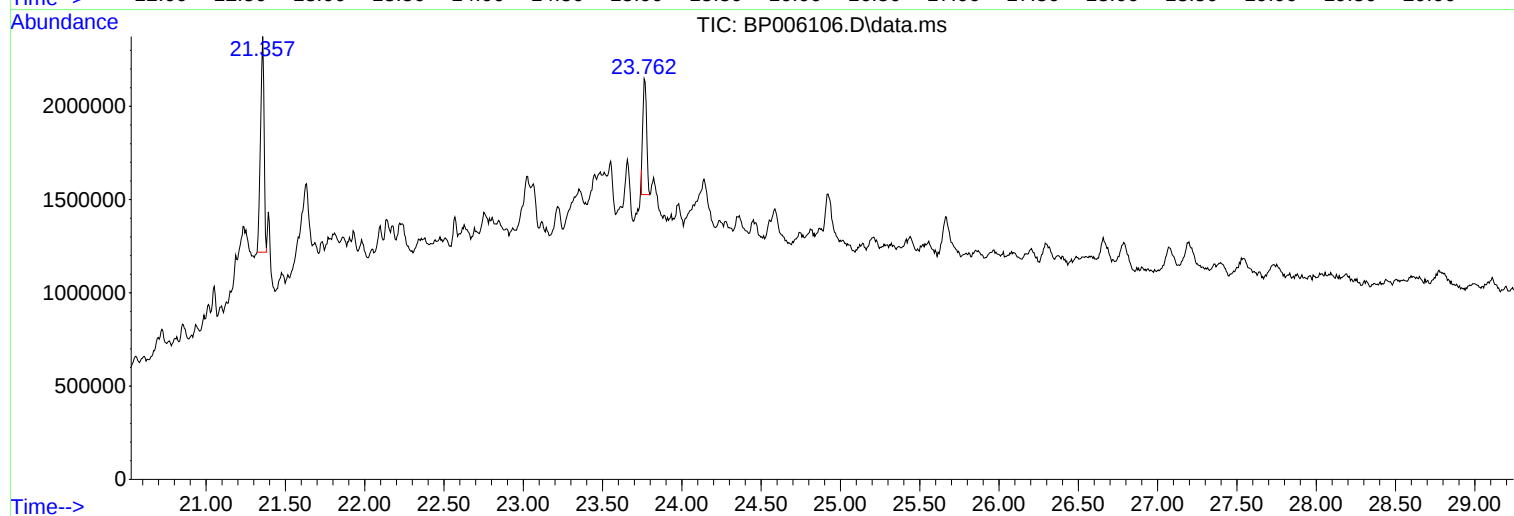
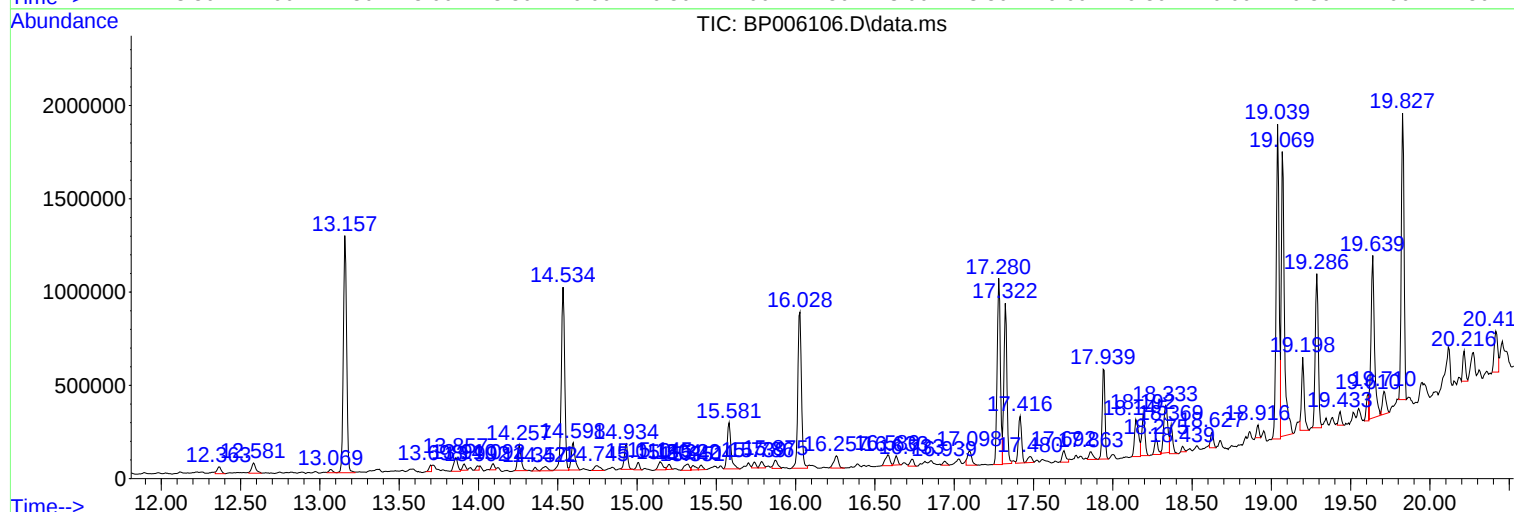
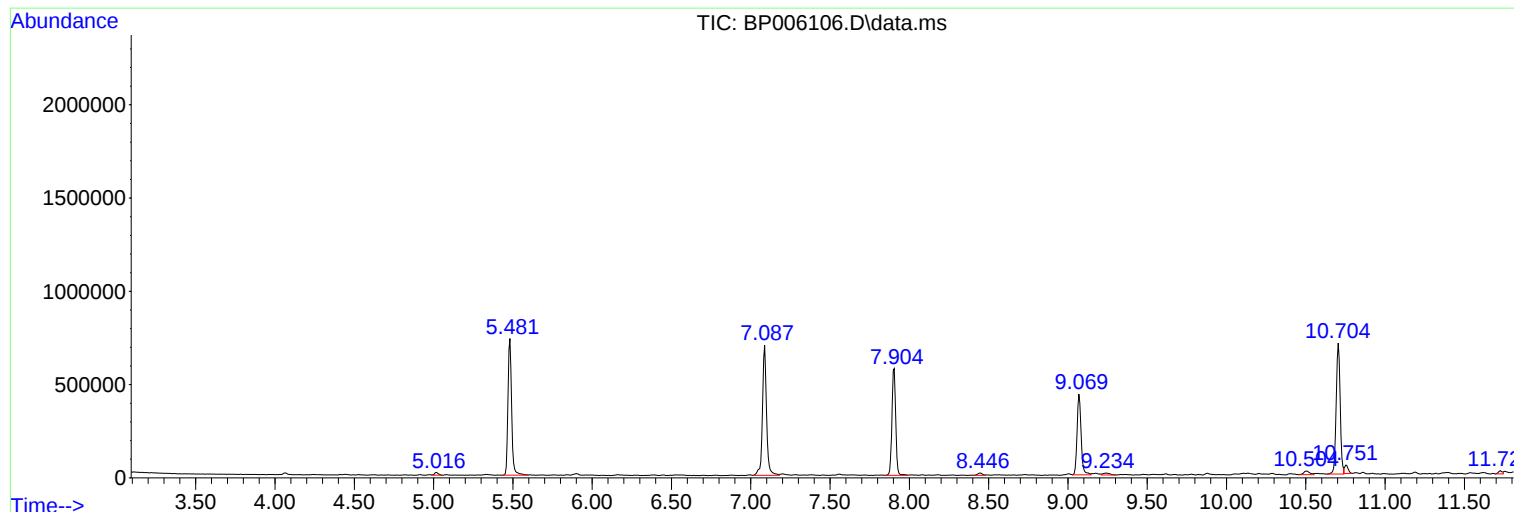
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

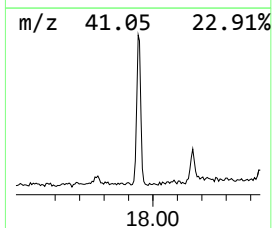
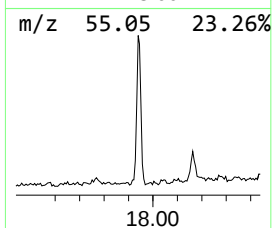
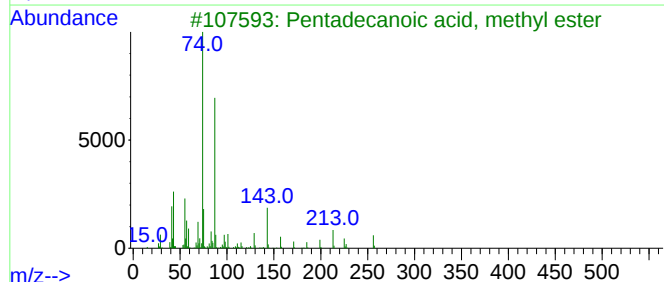
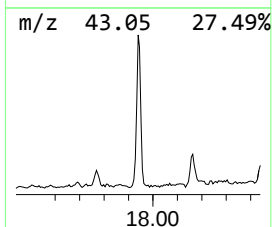
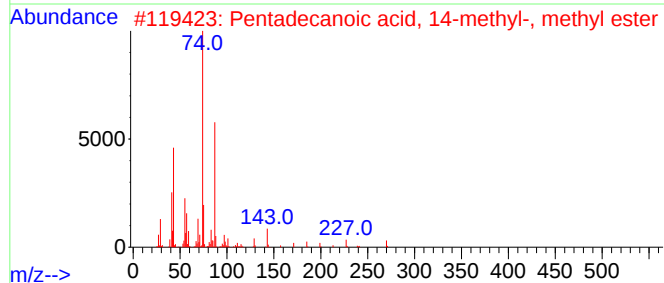
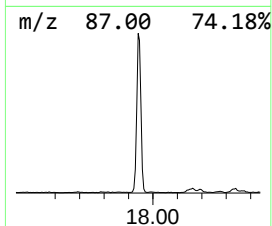
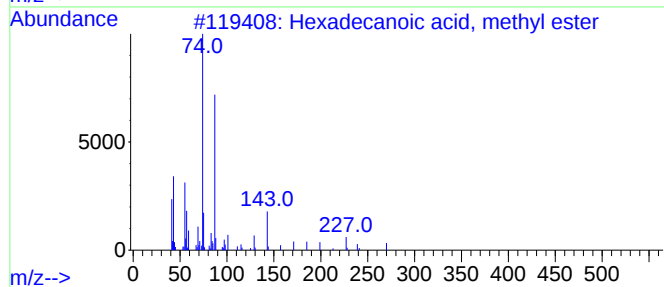
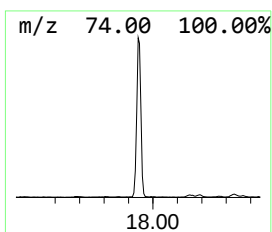
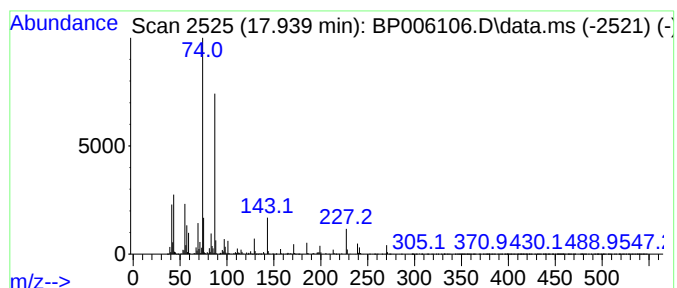
Instrument :
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TR-04-062821

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.939	8.13 ng	583369	Phenanthrene-d10	17.280	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	92
3	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	83
5	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	83



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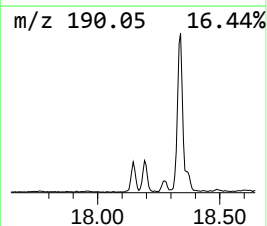
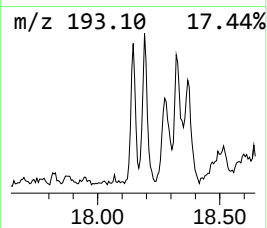
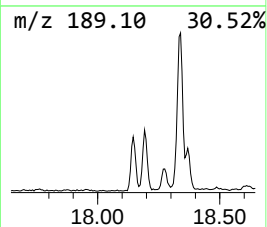
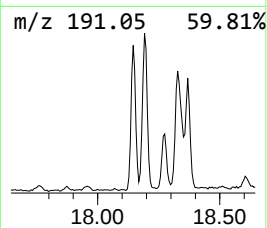
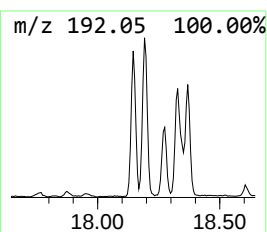
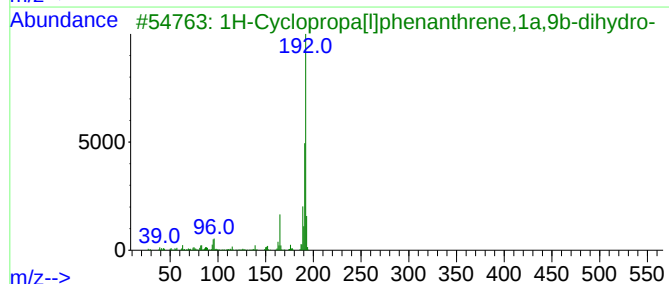
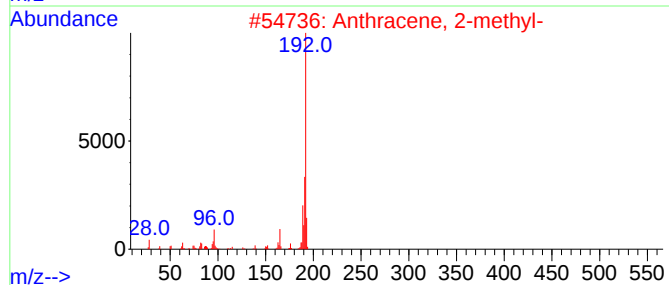
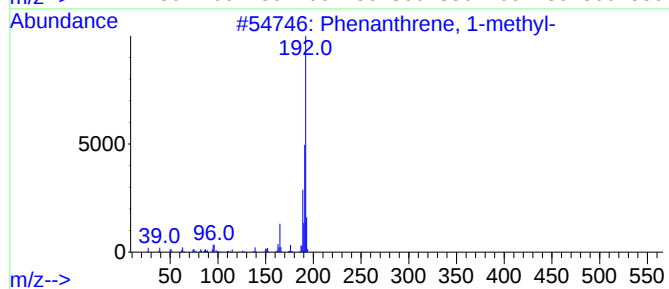
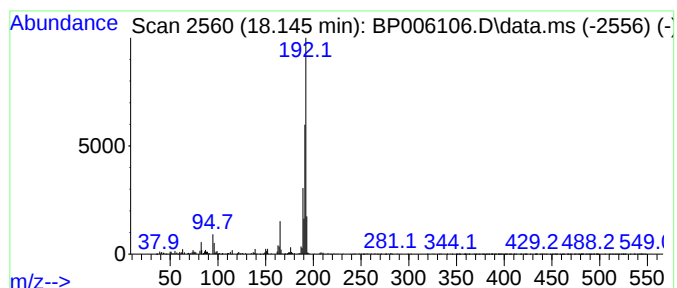
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.145	5.10 ng	365998	Phenanthrene-d10	17.280	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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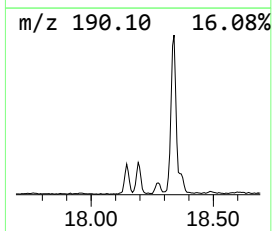
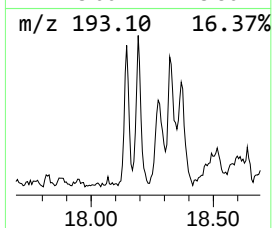
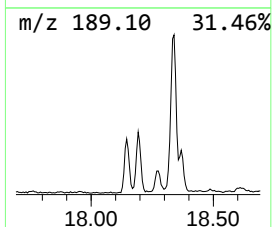
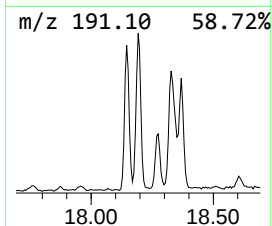
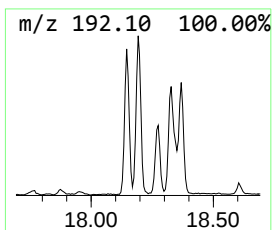
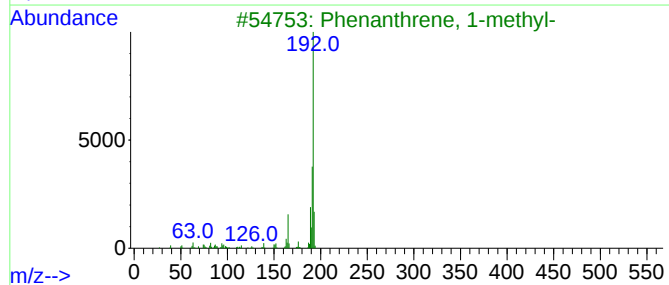
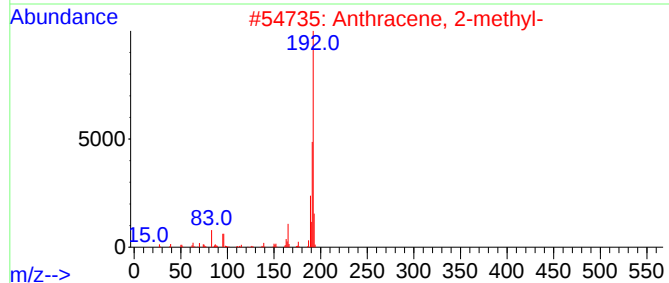
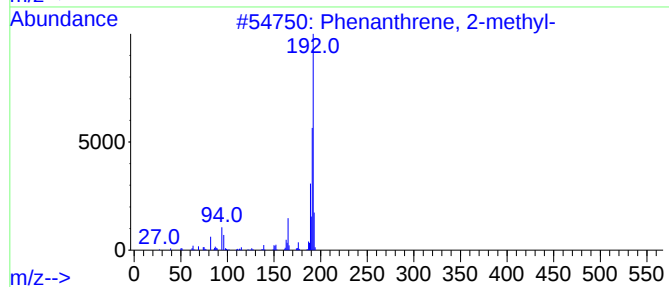
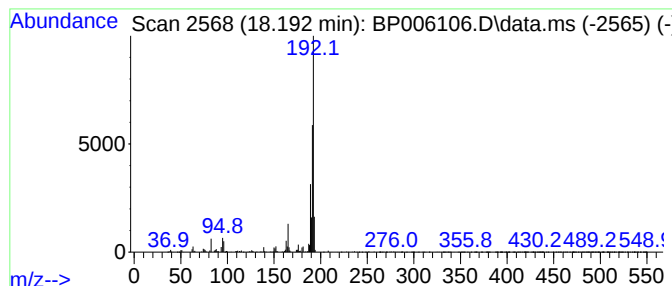
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	4.18 ng	300224	Phenanthrene-d10	17.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94



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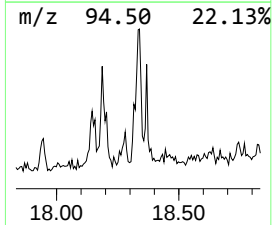
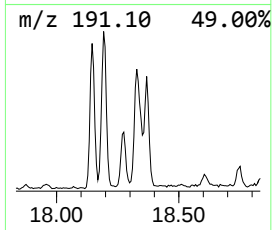
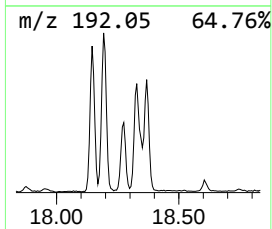
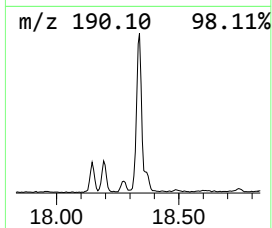
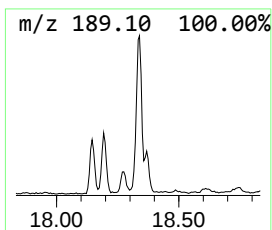
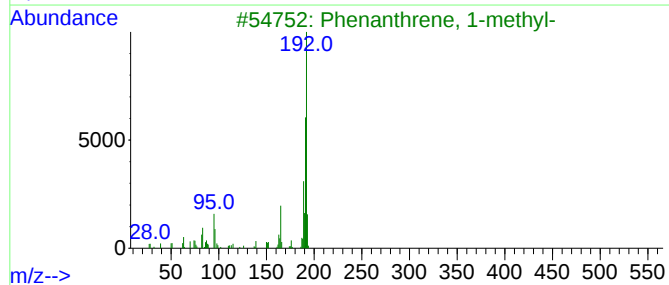
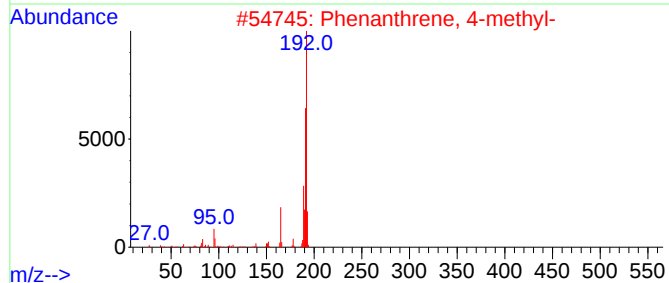
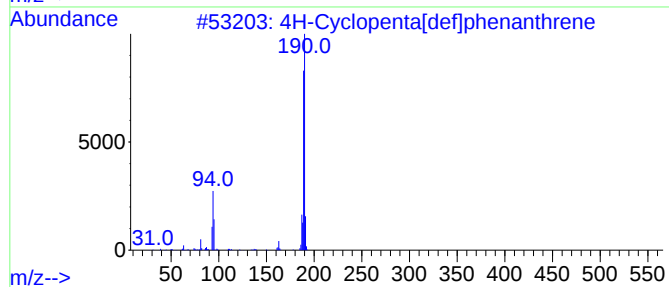
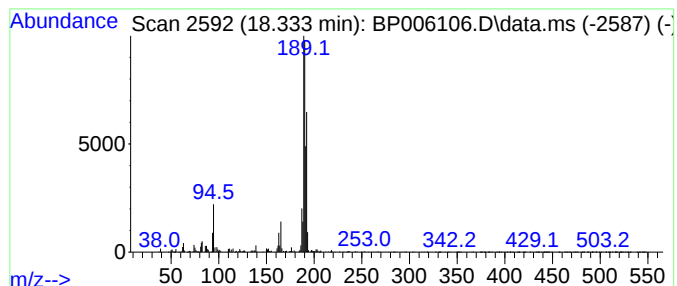
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.333	6.47 ng	463980	Phenanthrene-d10	17.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	35
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	30
5			Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TR-04-062821

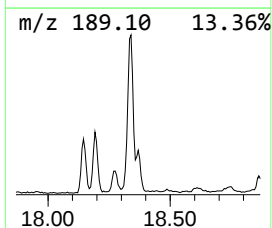
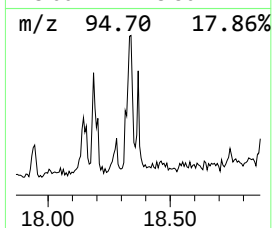
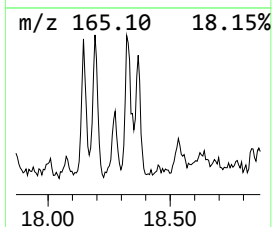
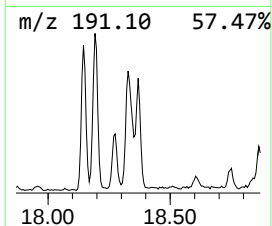
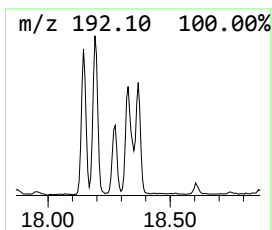
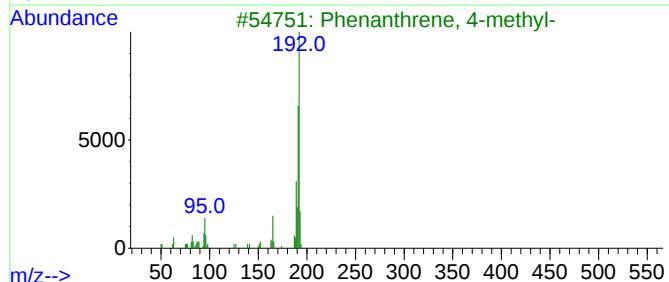
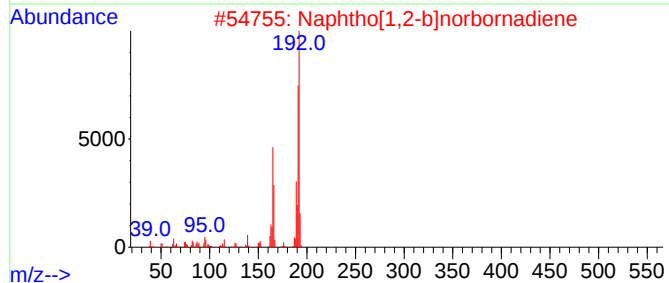
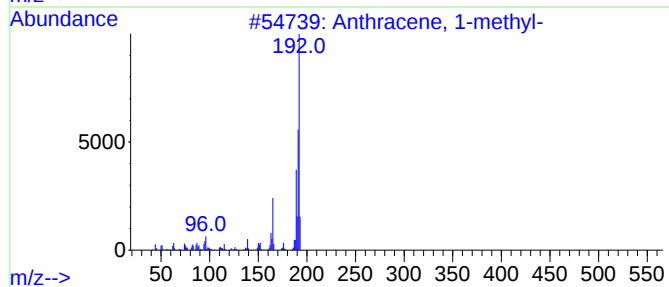
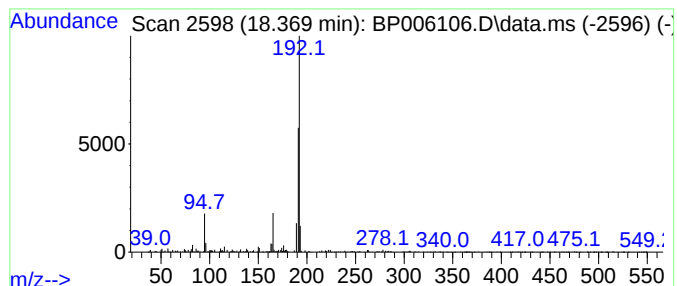
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	2.37 ng	170293	Phenanthrene-d10	17.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	72
2			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	72
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	72
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	68
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TR-04-062821

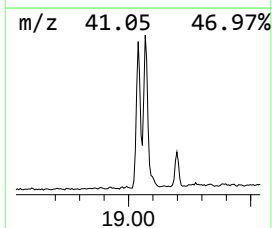
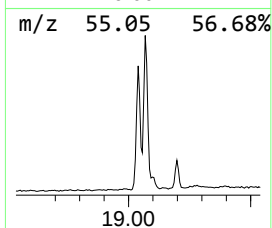
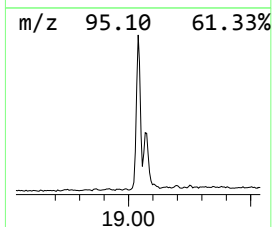
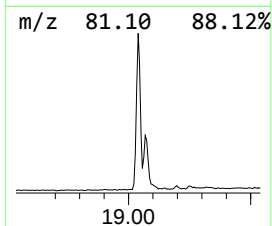
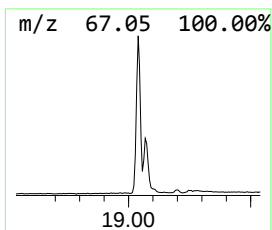
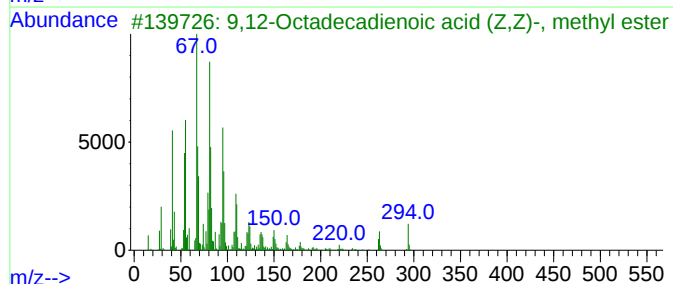
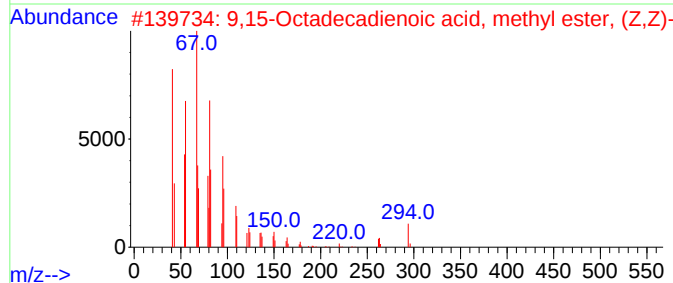
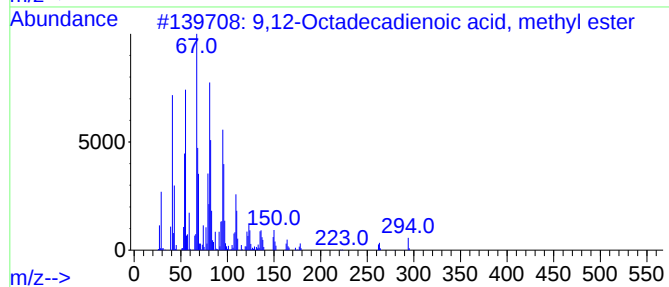
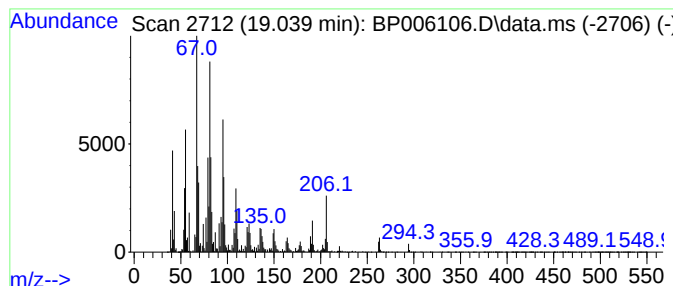
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.039	30.13 ng	2162070	Phenanthrene-d10	17.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
2			9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
5			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

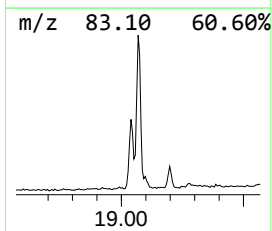
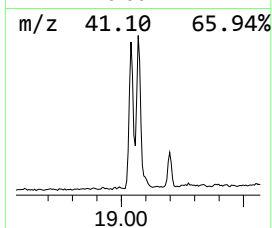
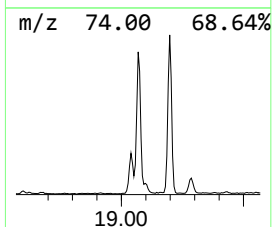
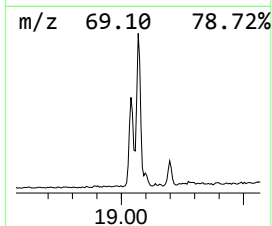
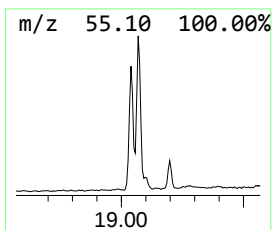
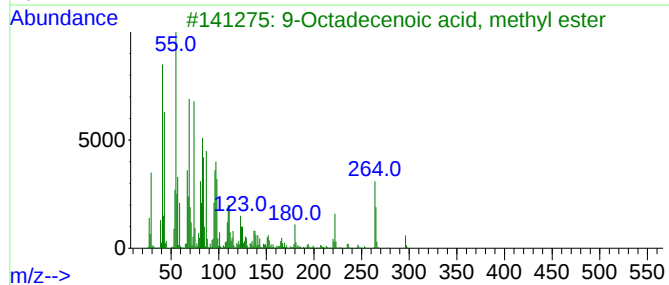
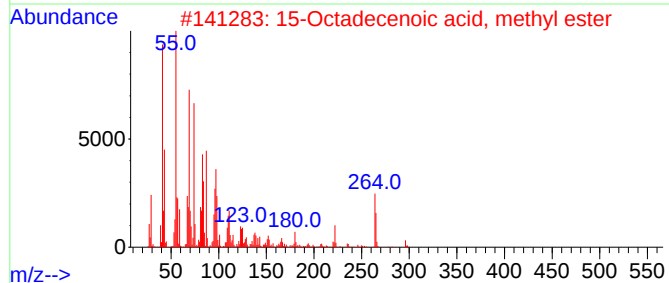
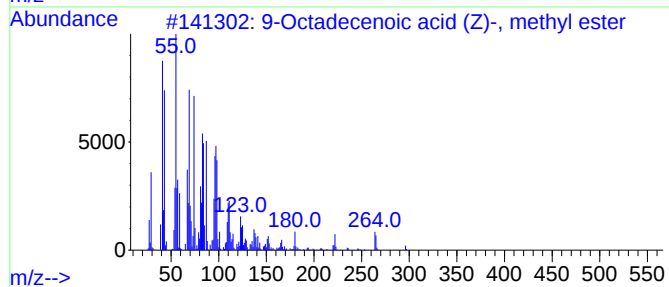
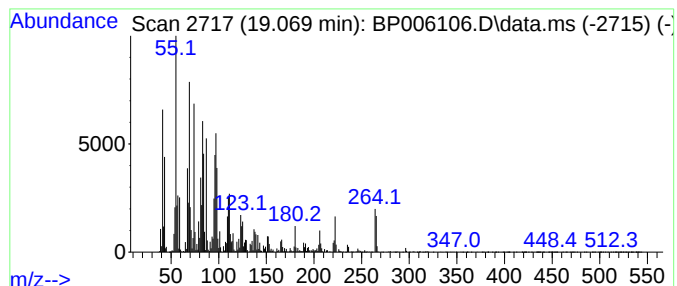
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.069	28.15 ng	2019990	Phenanthrene-d10	17.280	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
3	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
4	8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99
5	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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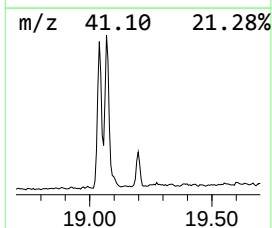
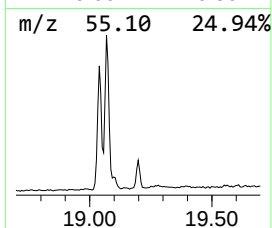
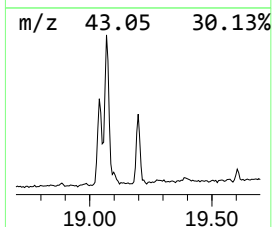
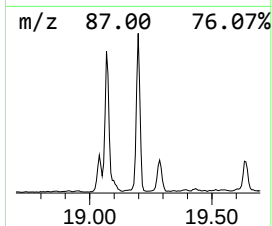
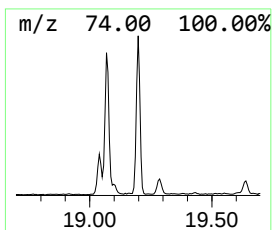
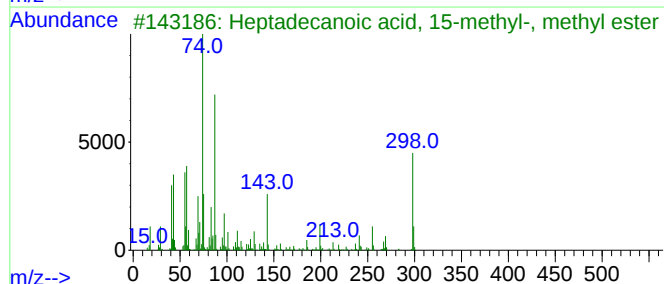
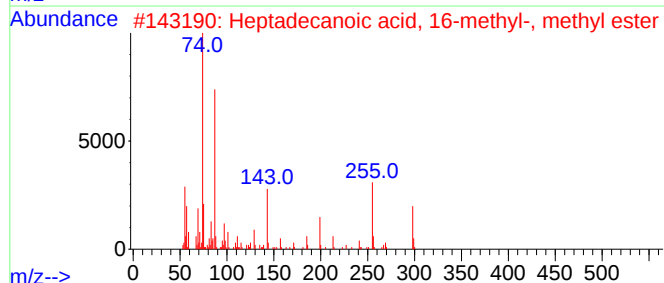
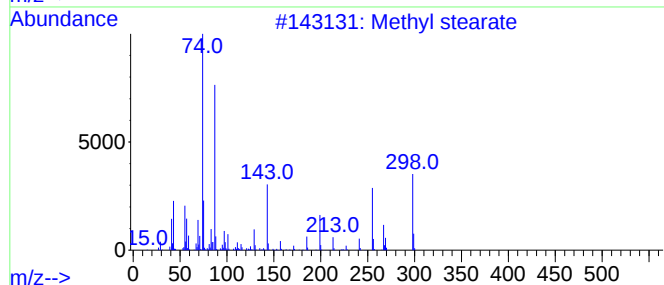
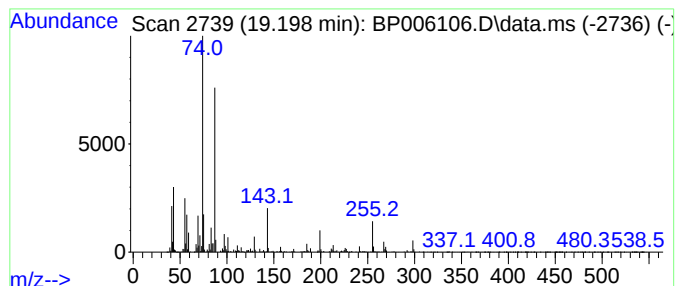
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.198	6.39 ng	458259	Phenanthrene-d10	17.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	93
3			Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	91
4			Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	91
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-04-062821

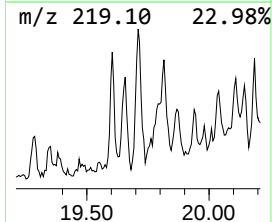
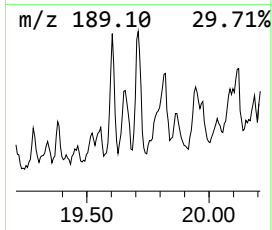
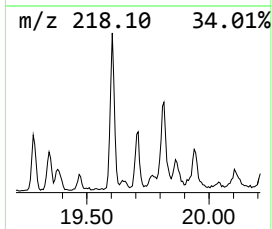
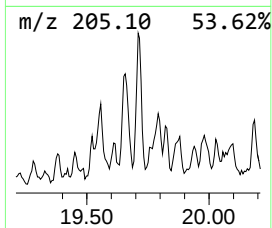
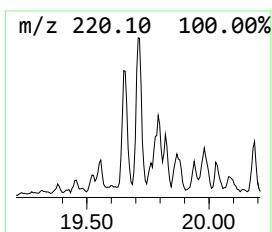
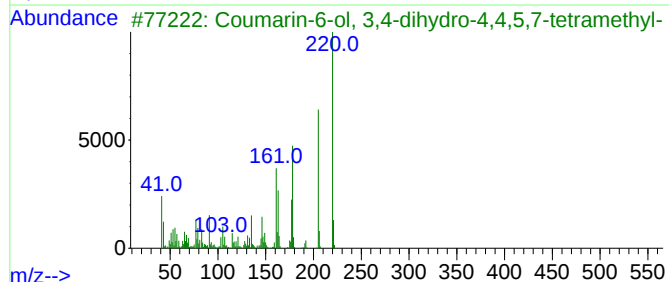
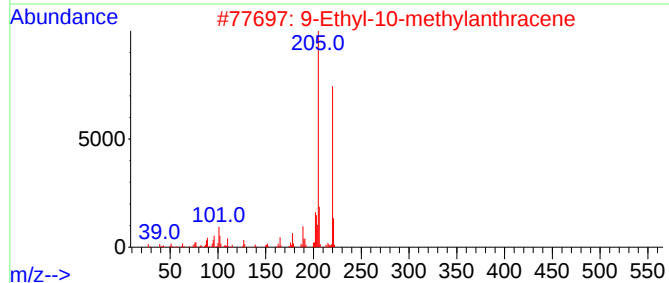
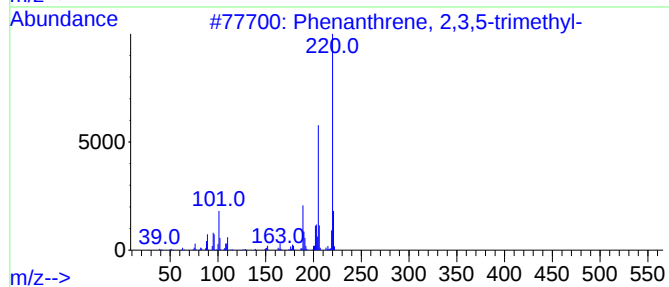
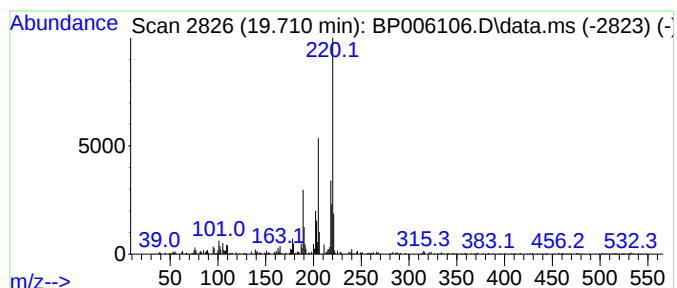
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.710	2.14 ng	191711	Chrysene-d12	21.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	83
2		9-Ethyl-10-methylantracene	220	C17H16	019713-49-6	50
3		Coumarin-6-ol, 3,4-dihydro-4,4,5...	220	C13H16O3	050442-68-7	49
4		Pyrrolo[1,2-a]thieno[2,3-d]pyrim...	220	C11H12N2OS	056929-61-4	49
5		1-Penten-3-one, 4-methyl-1-[2,6...	220	C15H24O	1000132-20-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-04-062821

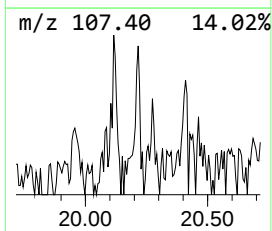
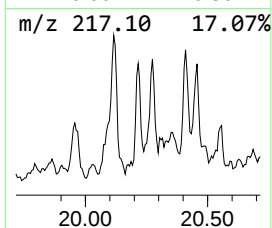
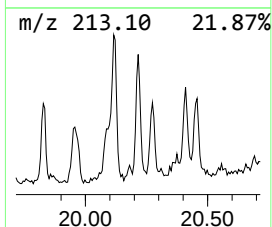
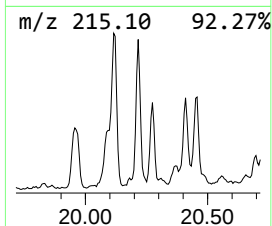
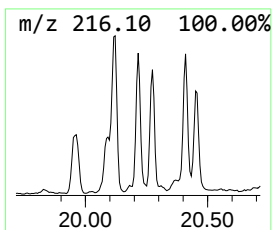
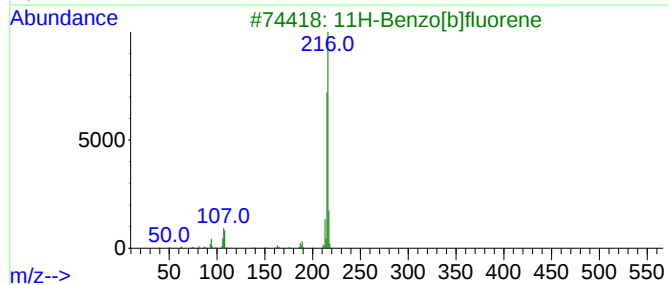
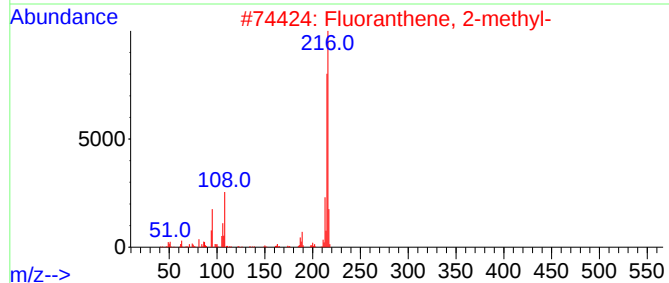
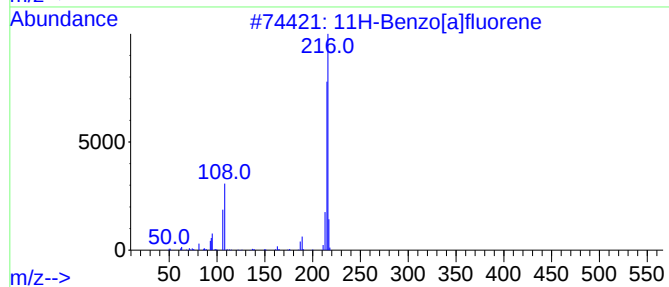
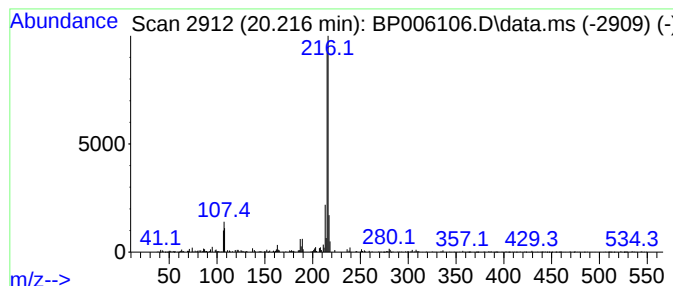
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.216	2.08 ng	186227	Chrysene-d12	21.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
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Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TR-04-062821

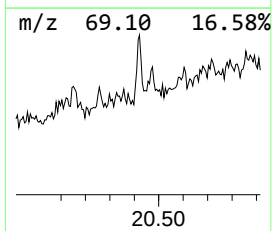
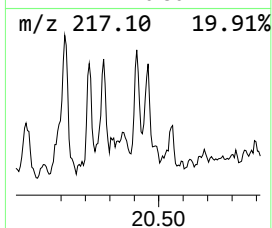
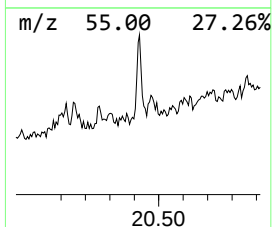
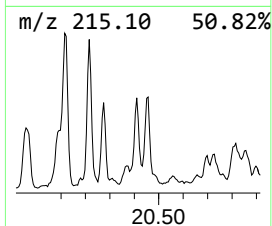
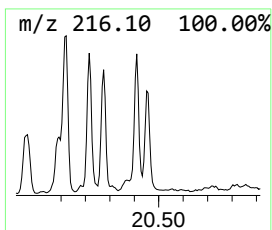
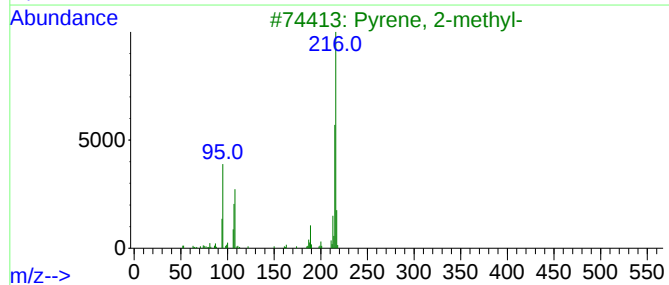
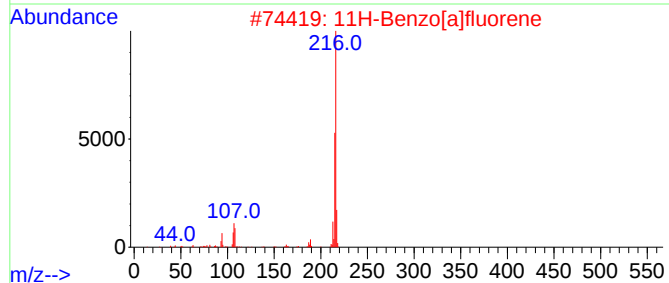
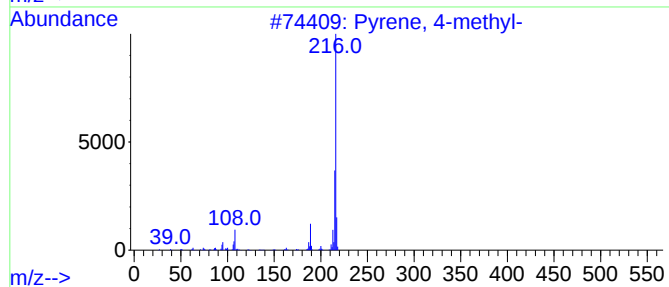
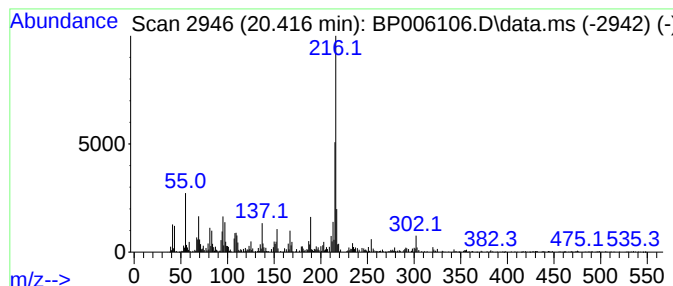
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.416	3.80 ng	340991	Chrysene-d12	21.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 4-methyl-	216	C17H12	003353-12-6	94
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	87
4			4-Hydroxy-6-methyl-1-pyridin-3-y...	216	C12H12N2O2	1000318-25-3	72
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
Data File : BP006106.D
Acq On : 29 Jun 2021 23:44
Operator : CG/JU
Sample : M2891-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TR-04-062821

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.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
Hexadecanoic ac...	17.939	8.1	ng	583369	4	17.280	1435200	20.0
Phenanthrene, 1...	18.145	5.1	ng	365998	4	17.280	1435200	20.0
Phenanthrene, 2...	18.192	4.2	ng	300224	4	17.280	1435200	20.0
4H-Cyclopenta[d...	18.333	6.5	ng	463980	4	17.280	1435200	20.0
Anthracene, 1-m...	18.369	2.4	ng	170293	4	17.280	1435200	20.0
9,12-Octadecadi...	19.039	30.1	ng	2162070	4	17.280	1435200	20.0
9-Octadecenoic ...	19.069	28.1	ng	2019990	4	17.280	1435200	20.0
Methyl stearate	19.198	6.4	ng	458259	4	17.280	1435200	20.0
Phenanthrene, 2...	19.710	2.1	ng	191711	5	21.357	1794470	20.0
11H-Benzo[a]flu...	20.216	2.1	ng	186227	5	21.357	1794470	20.0
Pyrene, 4-methyl-	20.416	3.8	ng	340991	5	21.357	1794470	20.0