

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
 Data File : BF136269.D
 Acq On : 29 Nov 2023 01:54
 Operator : CG\JU
 Sample : 05507-01 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-11202023

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_F\Methods\8270-BF112723.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF136269.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.881	301	305	309	rBV	13158	17741	1.03%	0.113%
2	5.428	565	568	577	rVB	244283	206818	11.97%	1.313%
3	6.428	735	738	745	rBV	240621	197536	11.43%	1.254%
4	6.804	798	802	805	rBV	818576	626784	36.28%	3.979%
5	7.357	893	896	901	rVV	142867	111642	6.46%	0.709%
6	7.398	901	903	906	rVB	30257	21897	1.27%	0.139%
7	8.081	1014	1019	1022	rBV	880828	725275	41.98%	4.604%
8	8.104	1022	1023	1026	rVV	60002	30987	1.79%	0.197%
9	8.675	1117	1120	1124	rBV	39478	30185	1.75%	0.192%
10	8.792	1138	1140	1144	rVB	29030	22422	1.30%	0.142%
11	8.892	1152	1157	1161	rBV2	31124	33298	1.93%	0.211%
12	9.157	1198	1202	1205	rBV	281849	210506	12.18%	1.336%
13	9.245	1213	1217	1221	rBV2	52723	59102	3.42%	0.375%
14	9.498	1256	1260	1262	rVV2	24614	27721	1.60%	0.176%
15	9.563	1266	1271	1273	rVV4	17374	22175	1.28%	0.141%
16	9.775	1304	1307	1310	rVB	53722	38607	2.23%	0.245%
17	9.839	1314	1318	1321	rVV	706351	621604	35.98%	3.946%
18	9.869	1321	1323	1327	rVB	96957	72023	4.17%	0.457%
19	9.998	1340	1345	1349	rBV5	9134	17598	1.02%	0.112%
20	10.039	1349	1352	1357	rVV	74233	71408	4.13%	0.453%
21	10.275	1389	1392	1395	rVB	47638	37211	2.15%	0.236%
22	10.386	1408	1411	1414	rBV	127614	98574	5.71%	0.626%
23	10.545	1435	1438	1442	rVB	25030	24005	1.39%	0.152%
24	10.622	1448	1451	1456	rVB	135497	130058	7.53%	0.826%
25	10.751	1470	1473	1482	rVB2	54689	66455	3.85%	0.422%
26	10.939	1501	1505	1507	rBV2	18232	19680	1.14%	0.125%
27	10.998	1512	1515	1525	rBV3	113642	120868	7.00%	0.767%
28	11.198	1542	1549	1552	rBV3	55501	81856	4.74%	0.520%
29	11.228	1552	1554	1558	rVB2	39409	38465	2.23%	0.244%
30	11.328	1568	1571	1573	rBV	817383	663862	38.42%	4.214%
31	11.351	1573	1575	1579	rVV	762904	592256	34.28%	3.760%
32	11.404	1579	1584	1587	rVV	175176	150716	8.72%	0.957%
33	11.439	1587	1590	1594	rVV3	20178	20293	1.17%	0.129%
34	11.563	1605	1611	1613	rBV	71251	76313	4.42%	0.484%
35	11.633	1620	1623	1625	rBV	43044	35796	2.07%	0.227%
36	11.733	1637	1640	1645	rVB	247694	190355	11.02%	1.208%

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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_F\Methods\8270-BF112723.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.798	1647	1651	1653	rBV5	20572	30499	1.77%	0.194%
38	11.833	1654	1657	1659	rVV	78941	68039	3.94%	0.432%
39	11.880	1659	1665	1672	rVV2	1420070	1727766	100.00%	10.968%
40	11.945	1673	1676	1678	rVV	133286	144600	8.37%	0.918%

41	11.969	1678	1680	1684	rVB	43718	45623	2.64%	0.290%
42	12.039	1689	1692	1695	rVV	34492	31975	1.85%	0.203%
43	12.133	1705	1708	1710	rBV	45166	44532	2.58%	0.283%
44	12.151	1710	1711	1716	rVB3	29841	23031	1.33%	0.146%
45	12.280	1728	1733	1736	rBV6	25974	40611	2.35%	0.258%

46	12.386	1748	1751	1754	rBV	39080	43251	2.50%	0.275%
47	12.422	1754	1757	1759	rBV	802339	713203	41.28%	4.528%
48	12.445	1759	1761	1764	rVB	536628	394532	22.83%	2.505%
49	12.545	1773	1778	1780	rBV	838844	800847	46.35%	5.084%
50	12.569	1780	1782	1783	rVV	259813	265404	15.36%	1.685%

51	12.586	1783	1785	1795	rVV	962170	1218981	70.55%	7.738%
52	12.663	1795	1798	1802	rVV	434308	376334	21.78%	2.389%
53	12.775	1812	1817	1820	rBV	629108	645026	37.33%	4.095%
54	12.898	1835	1838	1840	rBV	45710	49420	2.86%	0.314%
55	12.922	1840	1842	1849	rVB	310291	305337	17.67%	1.938%

56	12.992	1850	1854	1858	rBV3	35724	64418	3.73%	0.409%
57	13.027	1858	1860	1863	rVB	23845	18523	1.07%	0.118%
58	13.110	1871	1874	1877	rVB	112393	108672	6.29%	0.690%
59	13.174	1881	1885	1888	rBV2	121579	121289	7.02%	0.770%
60	13.210	1888	1891	1894	rBV	61889	60724	3.51%	0.385%

61	13.304	1905	1907	1909	rBV	28728	25237	1.46%	0.160%
62	13.333	1910	1912	1916	rVB2	37770	33960	1.97%	0.216%
63	13.369	1916	1918	1922	rBV4	25731	29132	1.69%	0.185%
64	13.410	1922	1925	1930	rBV	64369	67021	3.88%	0.425%
65	13.516	1940	1943	1945	rBV2	33555	31907	1.85%	0.203%

66	13.727	1976	1979	1982	rBV2	61892	70019	4.05%	0.444%
67	13.757	1982	1984	1986	rBV	40314	27259	1.58%	0.173%
68	13.792	1986	1990	1993	rBV3	37761	59562	3.45%	0.378%
69	13.827	1994	1996	1997	rBV	38112	30066	1.74%	0.191%
70	13.980	2018	2022	2025	rBV2	1160794	1356335	78.50%	8.610%

71	14.010	2025	2027	2029	rVB	230994	181511	10.51%	1.152%
72	14.086	2038	2040	2043	rVB	47652	41597	2.41%	0.264%
73	14.627	2130	2132	2137	rVB	101184	99084	5.73%	0.629%
74	15.033	2197	2201	2204	rBV	138286	209586	12.13%	1.331%
75	15.339	2250	2253	2258	rVB	90690	120850	6.99%	0.767%

76	15.404	2261	2264	2268	rVB	90266	105275	6.09%	0.668%
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.468	2271	2275	2278	rBV	331444	409178	23.68%	2.598%
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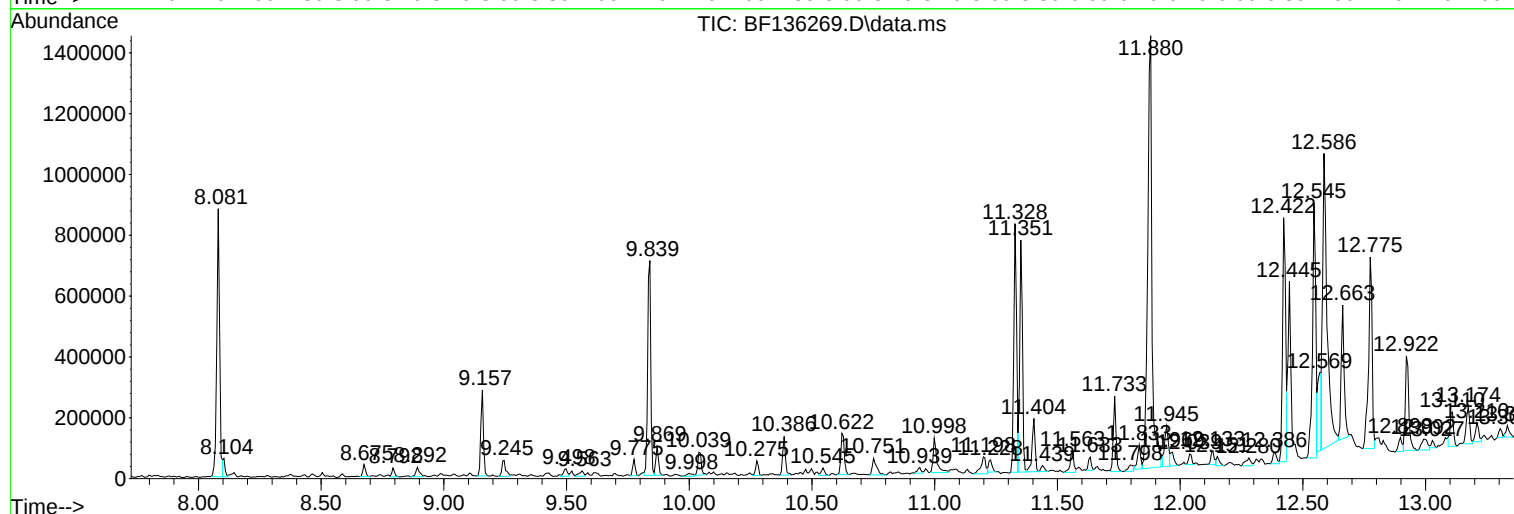
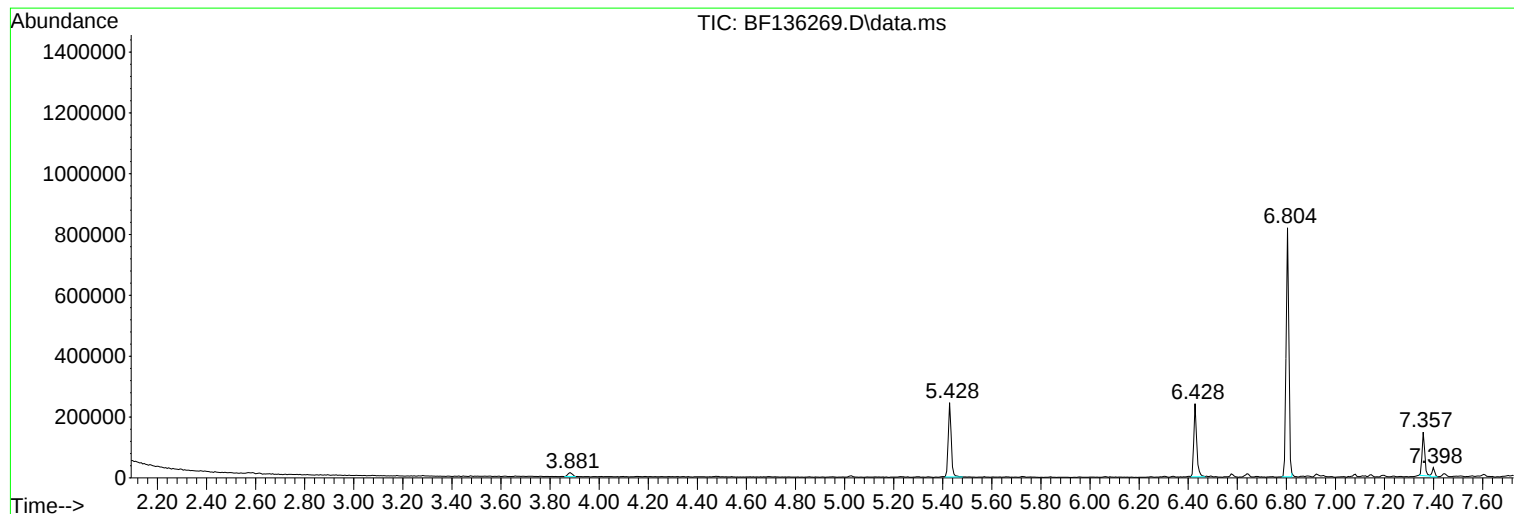
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Instrument :
BNA_F
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
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Instrument :
BNA_F
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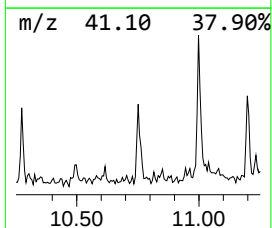
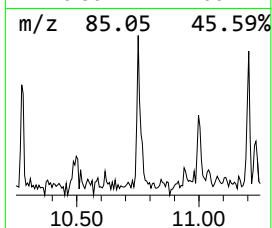
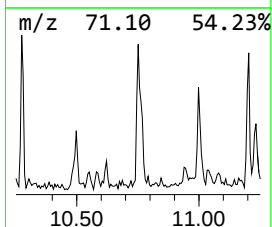
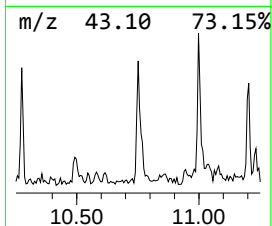
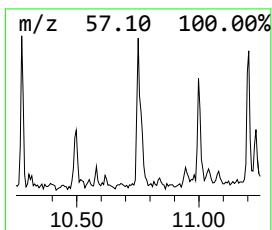
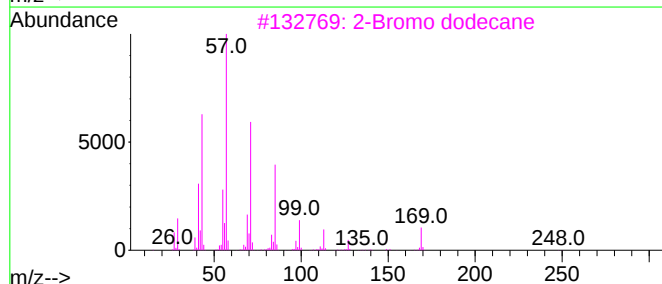
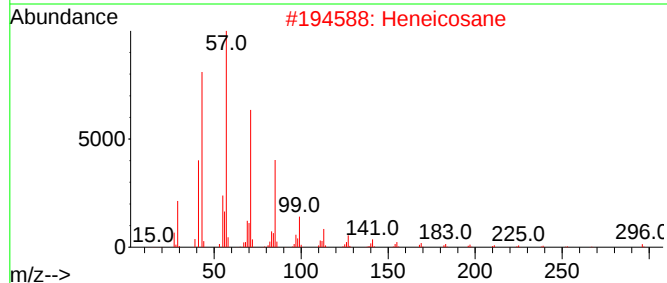
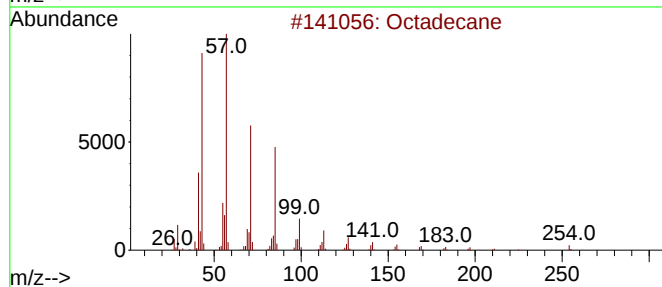
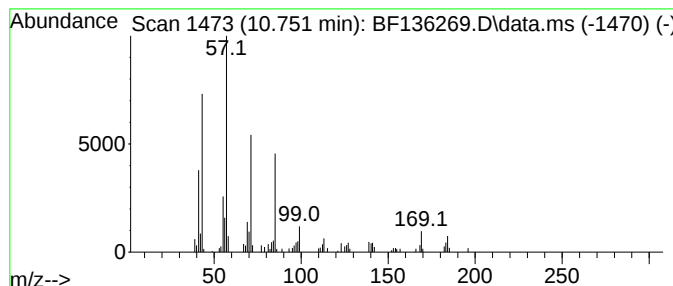
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.751	2.00 ng	66455	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	83
2			Heneicosane	296	C21H44	000629-94-7	83
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	80
4			Octacosane	394	C28H58	000630-02-4	80
5			Nonadecane	268	C19H40	000629-92-5	74



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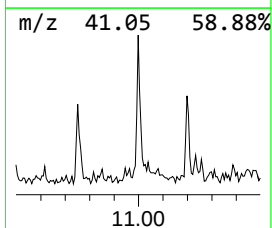
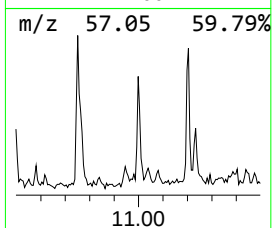
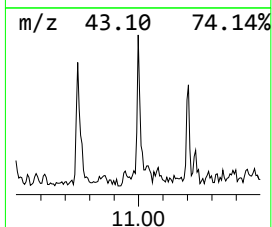
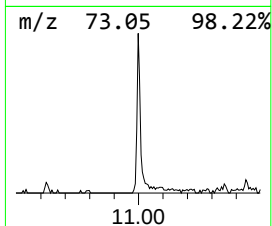
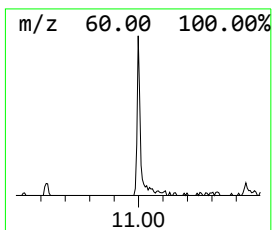
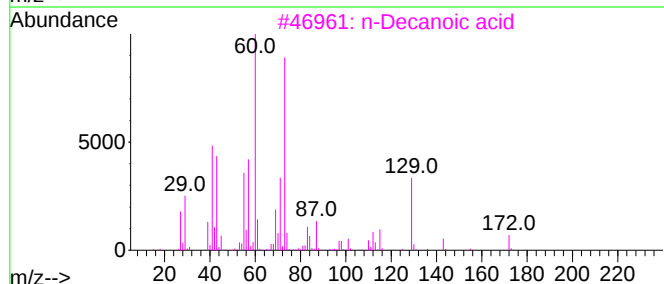
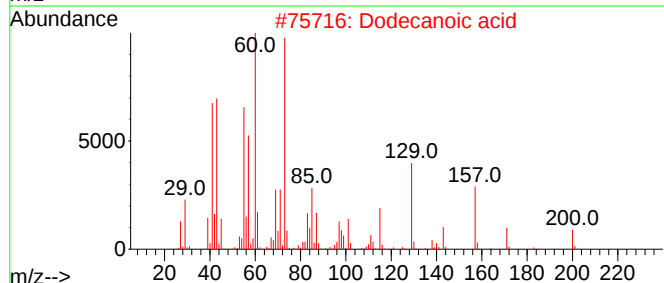
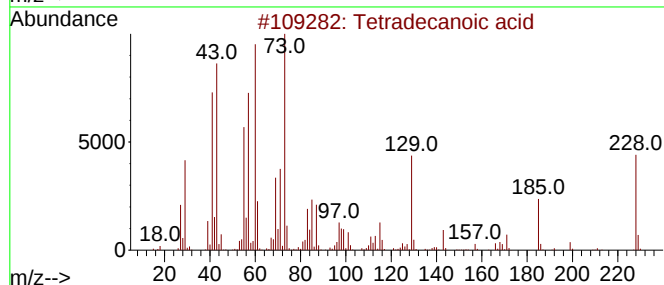
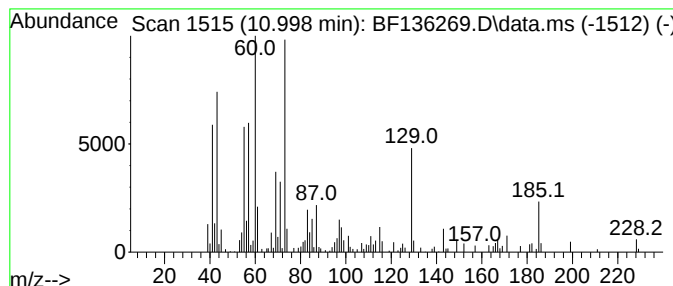
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tetradecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.998	3.64 ng	120868	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
2			Dodecanoic acid	200	C12H24O2	000143-07-7	72
3			n-Decanoic acid	172	C10H20O2	000334-48-5	64
4			Octanoic acid	144	C8H16O2	000124-07-2	43
5			1,2,3,4-Cyclopentanetetrol, (1.a...	134	C5H10O4	014003-71-5	35



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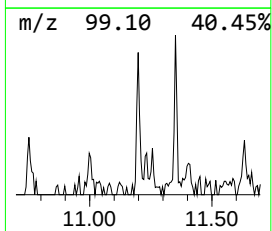
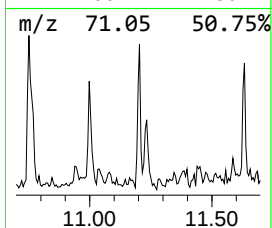
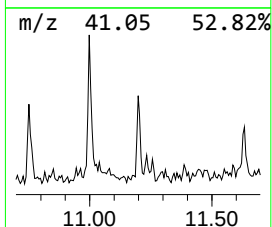
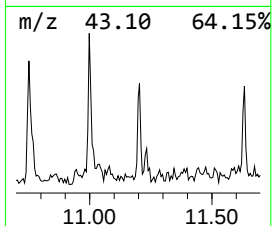
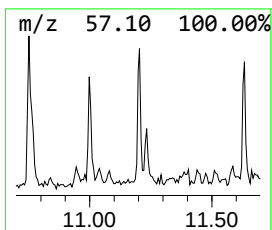
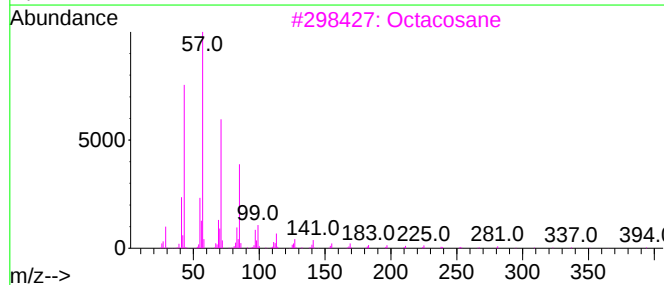
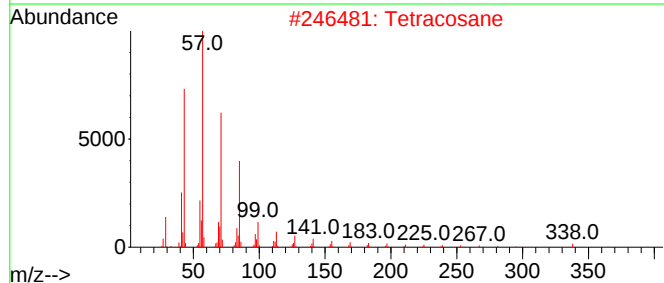
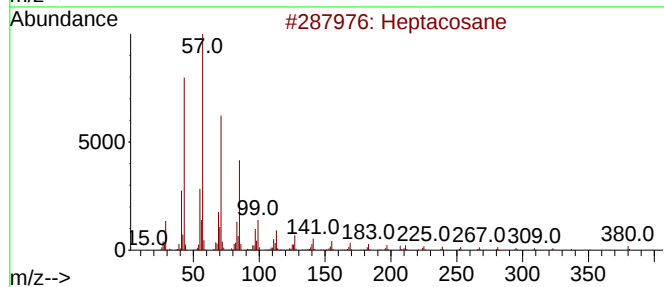
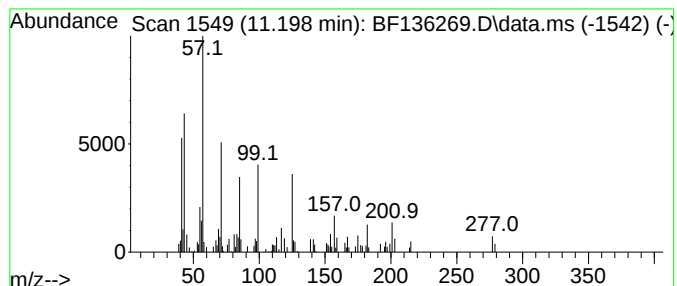
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown11.198 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	2.47 ng	81856	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	38
2			Tetracosane	338	C24H50	000646-31-1	38
3			Octacosane	394	C28H58	000630-02-4	38
4			Pentacosane	352	C25H52	000629-99-2	38
5			Heneicosane	296	C21H44	000629-94-7	38



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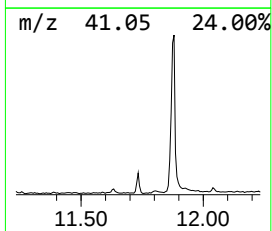
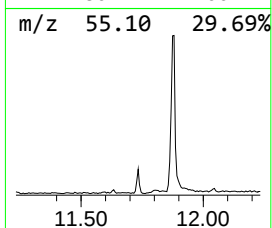
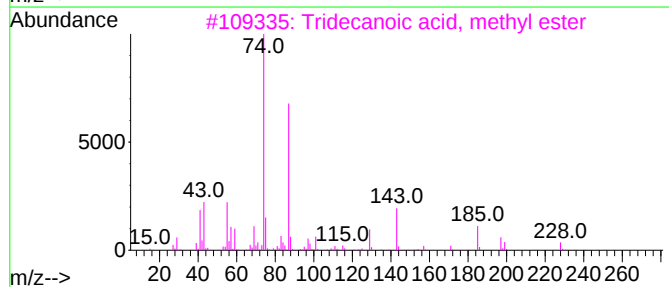
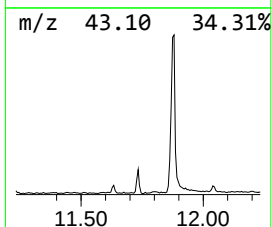
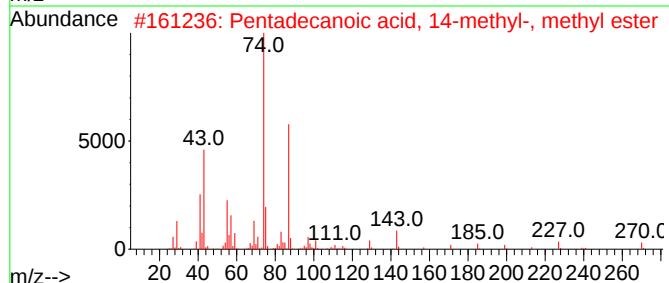
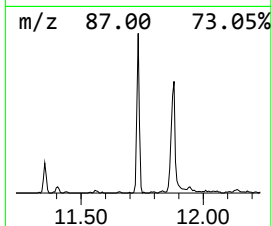
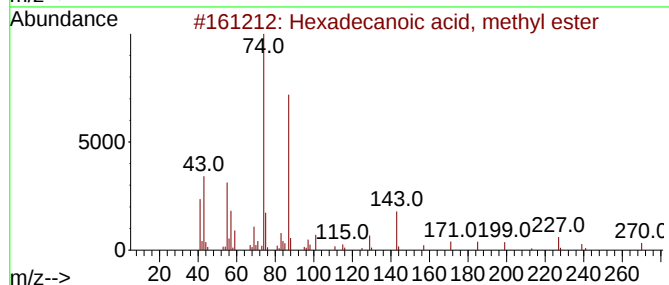
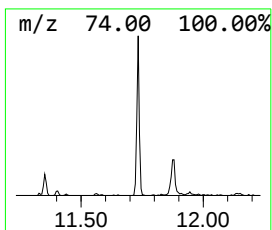
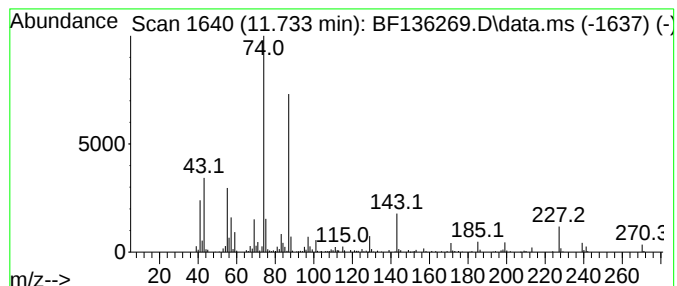
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.733	5.73 ng	190355	Phenanthrene-d10	11.328

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	93
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	83
4			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	81
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
Acq On : 29 Nov 2023 01:54
Operator : CG\JU
Sample : 05507-01 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SU-04-11202023

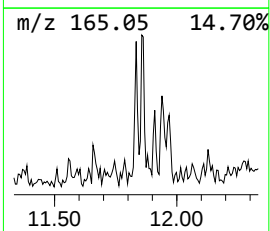
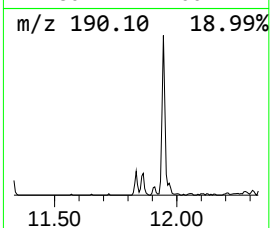
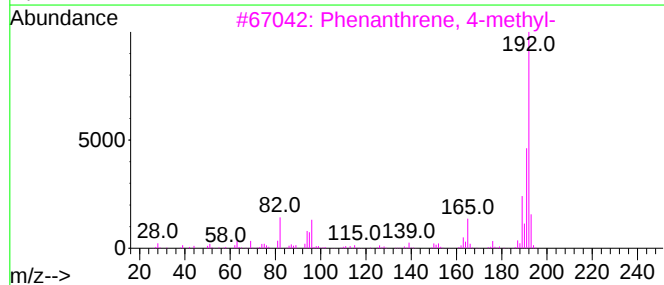
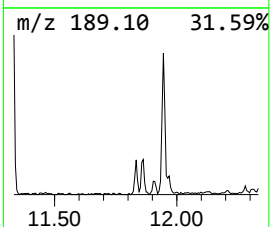
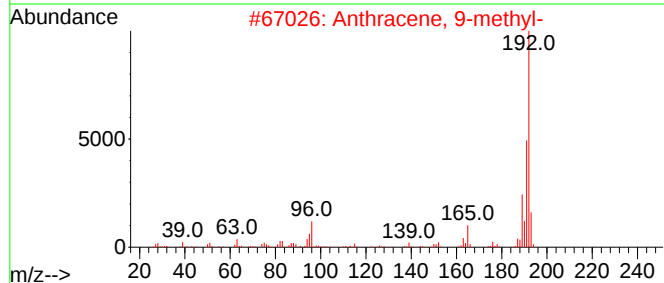
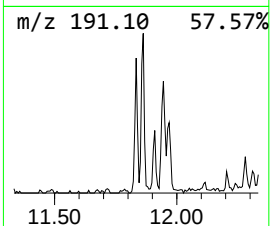
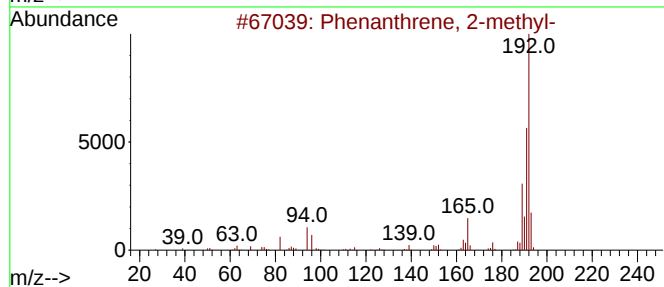
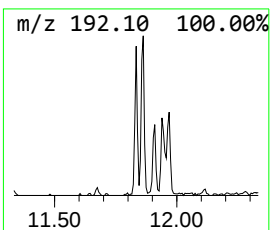
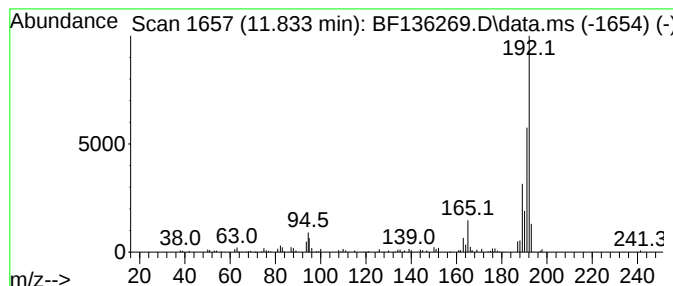
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.833	2.05 ng	68039	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	87
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
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Operator : CG\JU
Sample : 05507-01 10X
Misc :
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ClientSampleId :
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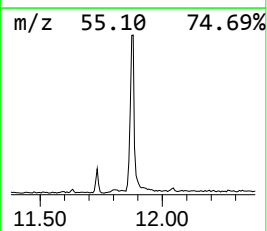
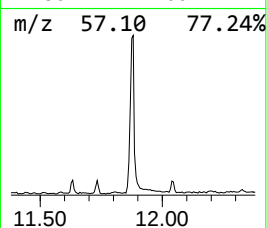
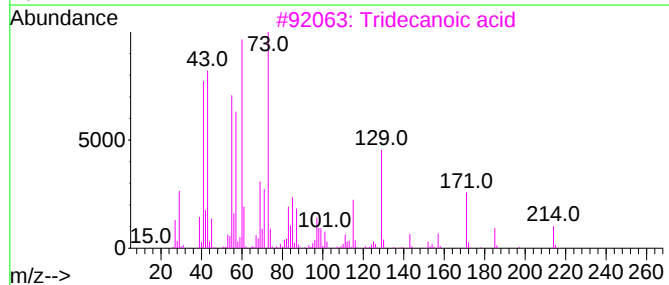
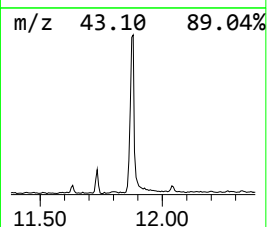
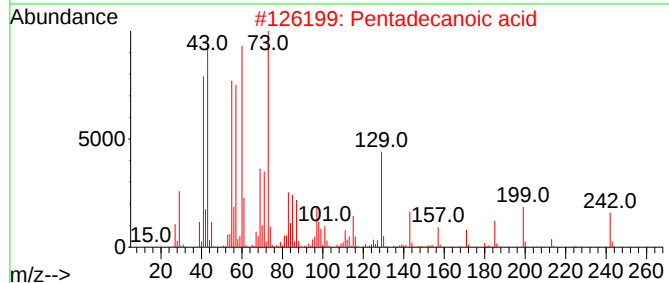
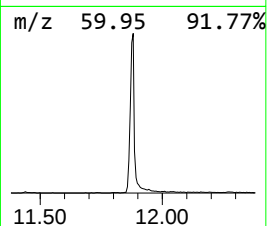
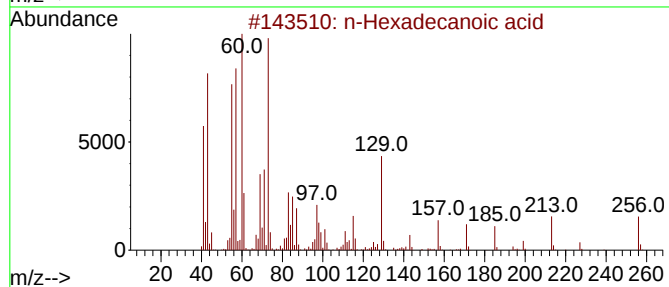
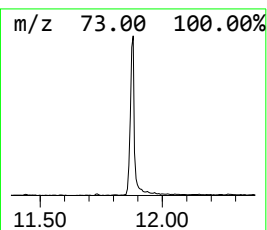
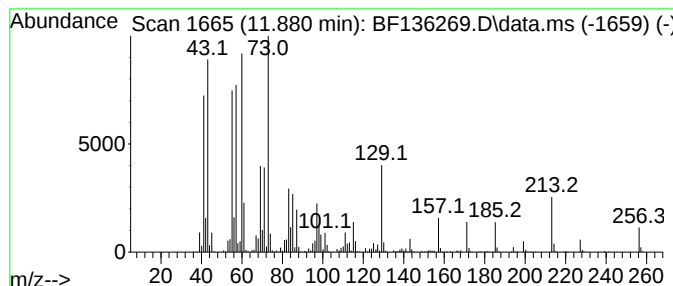
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	52.05 ng	1727770	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3			Tridecanoic acid	214	C13H26O2	000638-53-9	90
4			n-Decanoic acid	172	C10H20O2	000334-48-5	68
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
Acq On : 29 Nov 2023 01:54
Operator : CG\JU
Sample : 05507-01 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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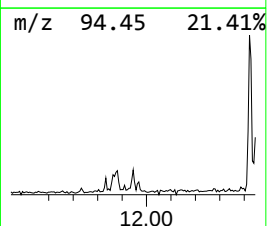
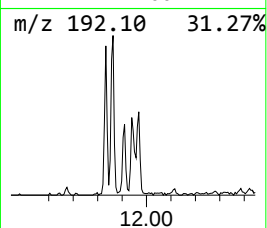
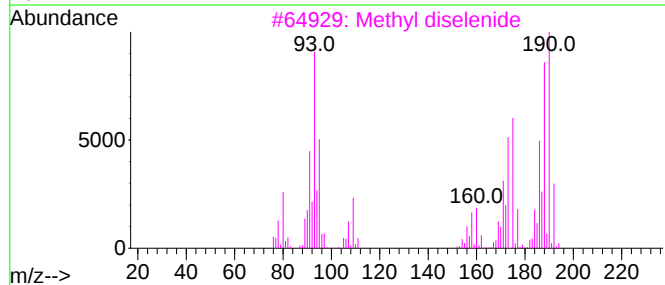
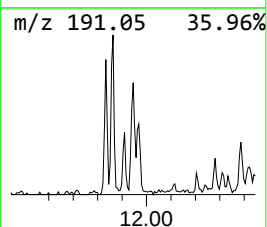
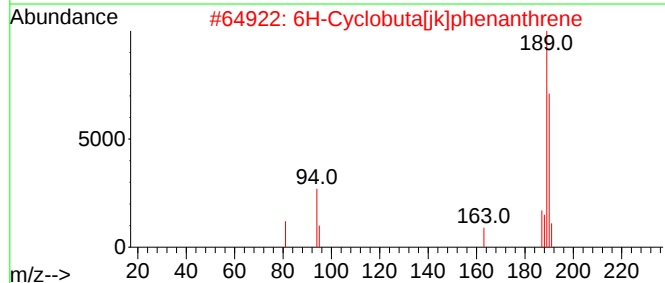
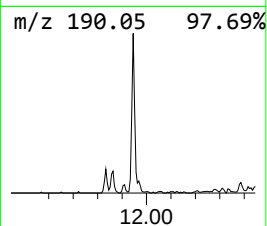
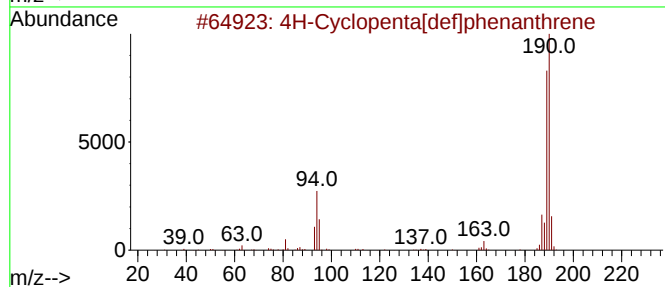
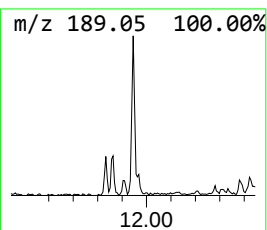
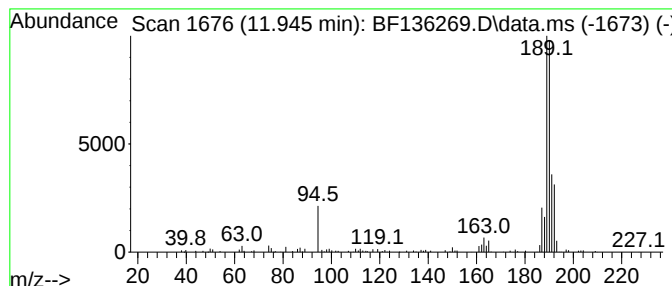
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	4.36 ng	144600	Phenanthrene-d10	11.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3			Methyl diselenide	190	C2H6Se2	007101-31-7	55
4			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
5			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
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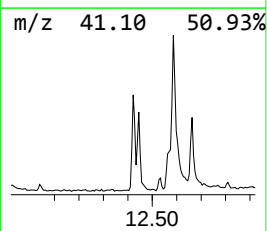
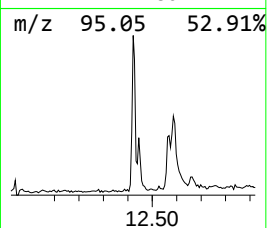
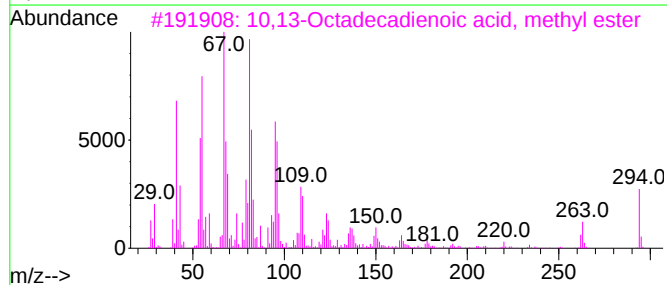
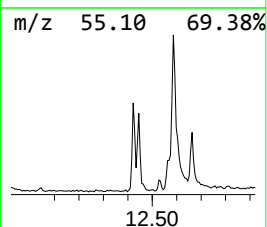
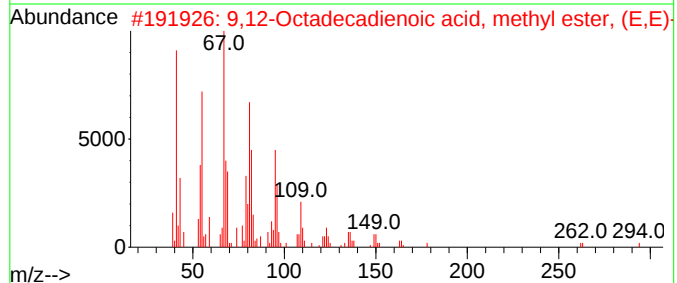
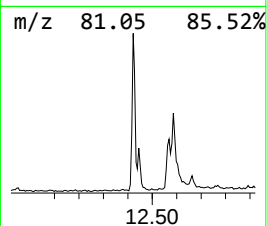
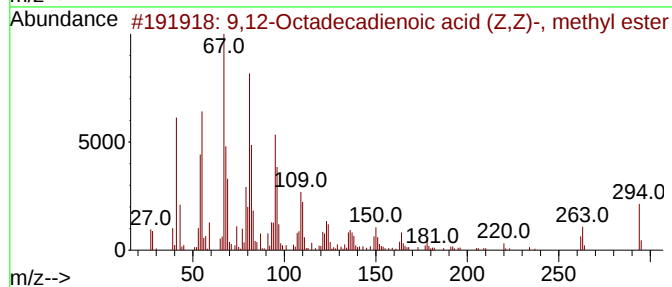
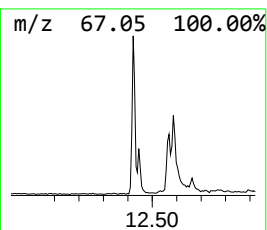
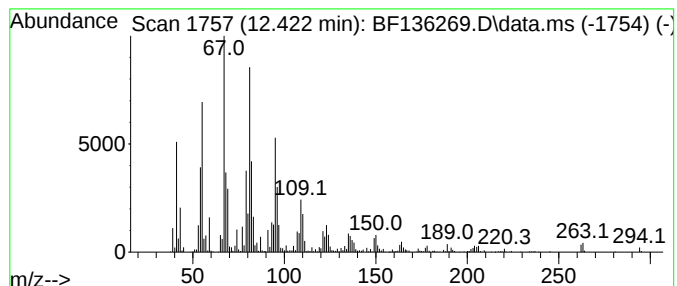
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,12-Octadecadienoic acid (... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.422	21.49 ng	713203	Phenanthrene-d10	11.328	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
4	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
Acq On : 29 Nov 2023 01:54
Operator : CG\JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SU-04-11202023

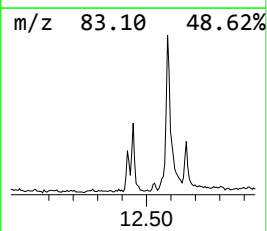
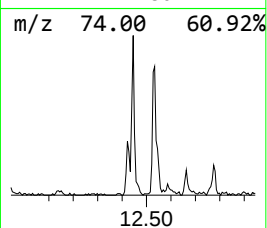
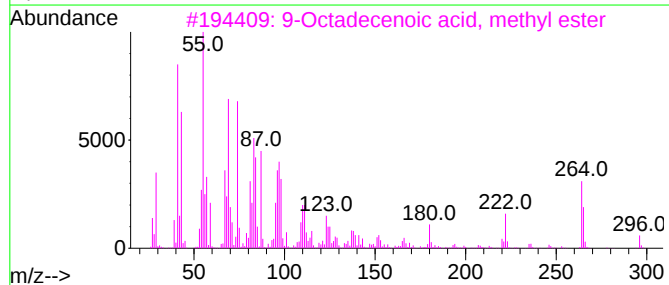
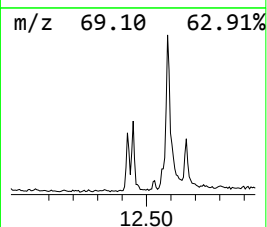
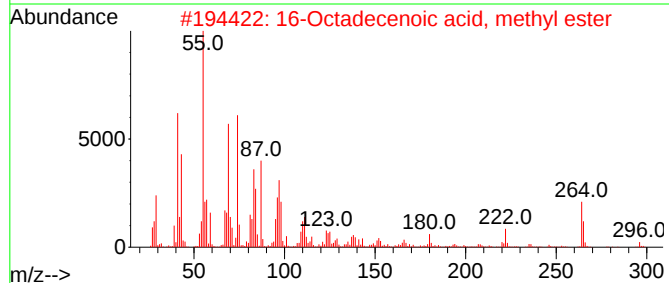
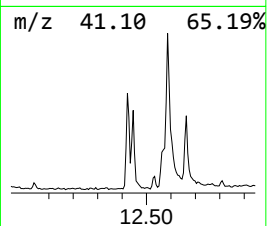
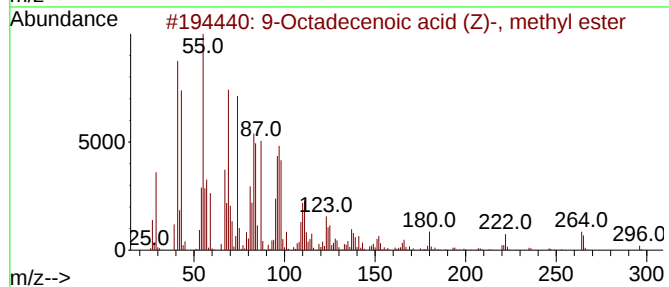
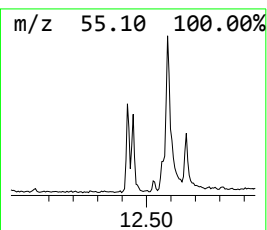
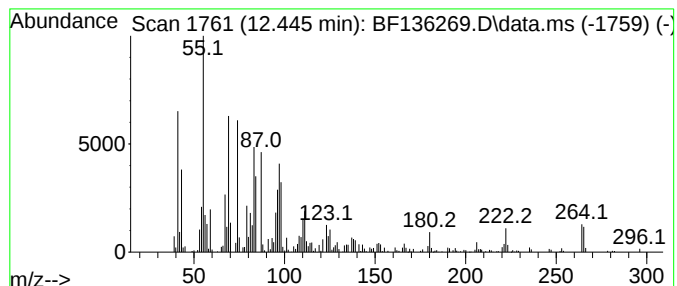
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	11.89 ng	394532	Phenanthrene-d10	11.328

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
3		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
4		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
5		11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
Acq On : 29 Nov 2023 01:54
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ALS Vial : 20 Sample Multiplier: 1

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SU-04-11202023

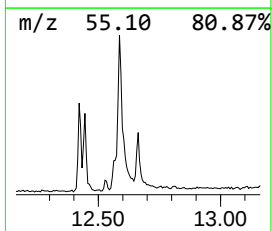
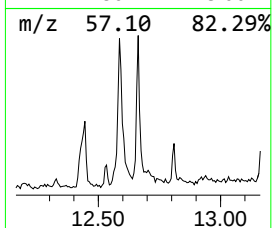
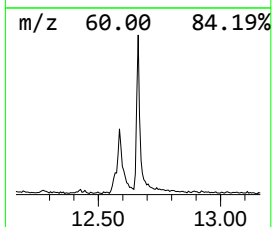
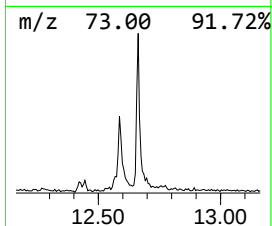
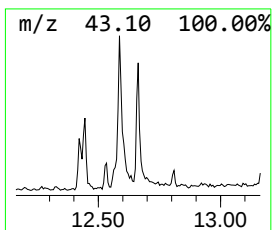
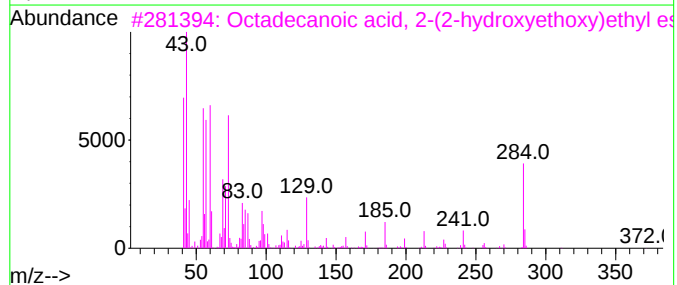
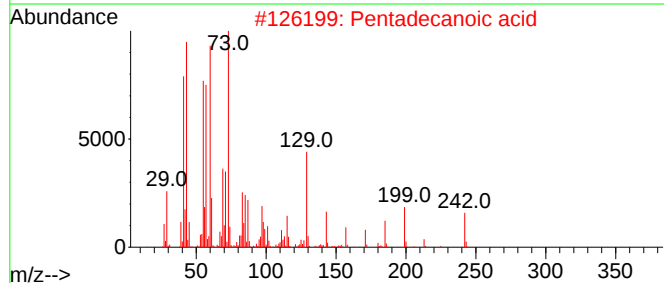
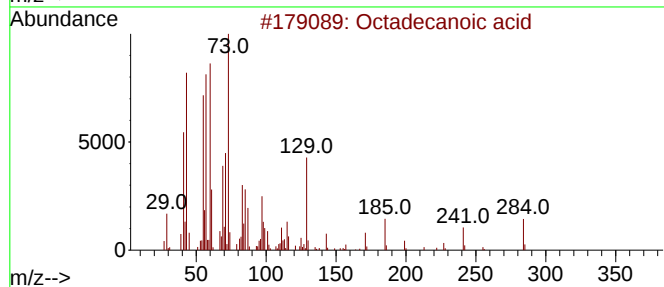
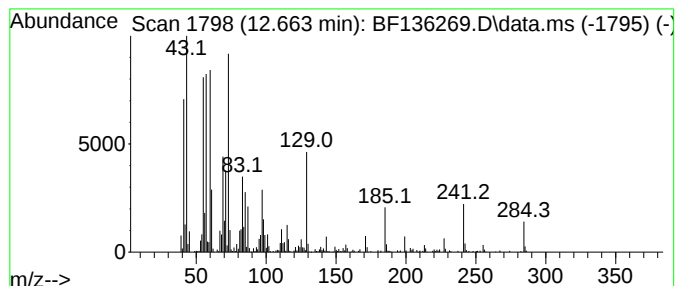
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.663	5.55 ng	376334	Chrysene-d12	13.980

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80
4			Tridecanoic acid	214	C13H26O2	000638-53-9	76
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
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ALS Vial : 20 Sample Multiplier: 1

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ClientSampleId :
SU-04-11202023

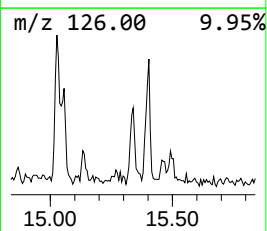
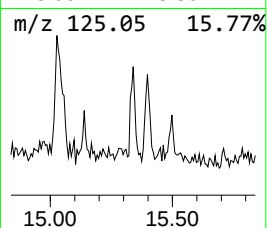
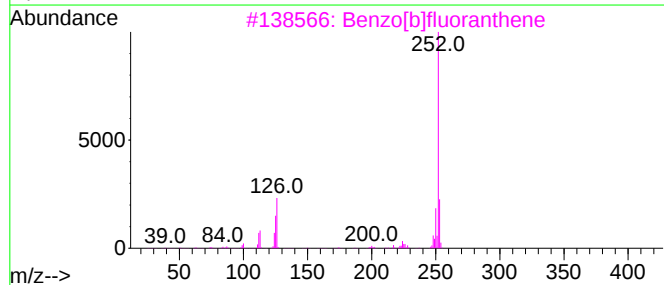
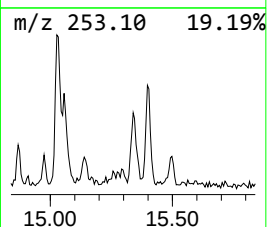
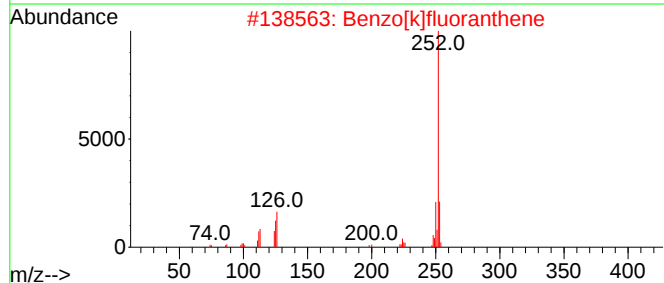
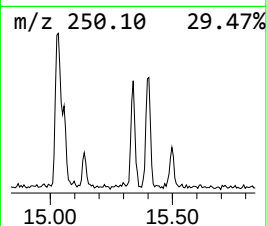
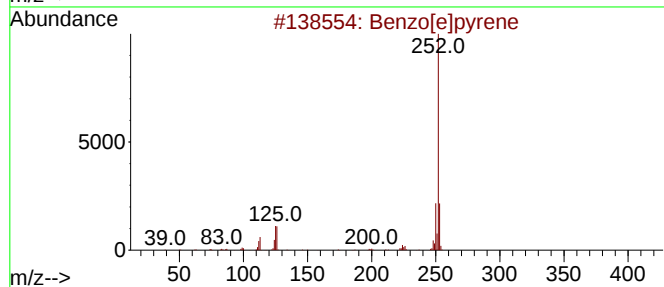
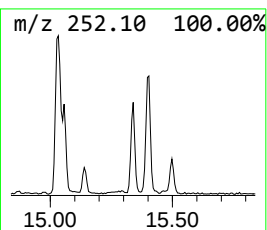
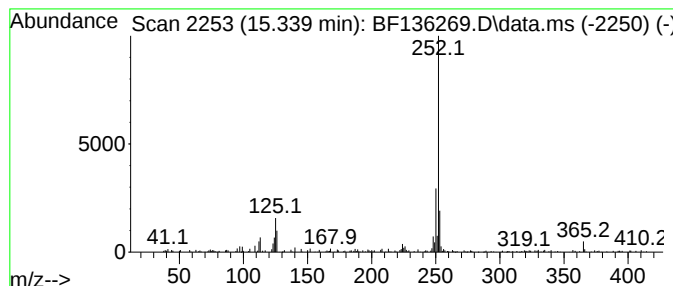
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.339	5.91 ng	120850	Perylene-d12	15.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	96
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Perylene	252	C20H12	000198-55-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
Data File : BF136269.D
Acq On : 29 Nov 2023 01:54
Operator : CG\JU
Sample : 05507-01 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-11202023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Octadecane	10.751	2.0	ng	66455	4	11.328	663862	20.0
Tetradecanoic acid	10.998	3.6	ng	120868	4	11.328	663862	20.0
unknown11.198	11.198	2.5	ng	81856	4	11.328	663862	20.0
Hexadecanoic ac...	11.733	5.7	ng	190355	4	11.328	663862	20.0
Phenanthrene, 2...	11.833	2.0	ng	68039	4	11.328	663862	20.0
n-Hexadecanoic ...	11.880	52.0	ng	1727770	4	11.328	663862	20.0
4H-Cyclopenta[d...	11.945	4.4	ng	144600	4	11.328	663862	20.0
9,12-Octadecadi...	12.422	21.5	ng	713203	4	11.328	663862	20.0
9-Octadecenoic ...	12.445	11.9	ng	394532	4	11.328	663862	20.0
Octadecanoic acid	12.663	5.5	ng	376334	5	13.980	1356340	20.0
Benzo[e]pyrene	15.339	5.9	ng	120850	6	15.468	409178	20.0