

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
 Data File : BF129837.D
 Acq On : 17 Aug 2022 10:51
 Operator : CG\JU
 Sample : N4199-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-02-081222

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-BF081422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129837.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.051	466	470	481	rVB	158861	205652	5.45%	0.676%
2	5.469	537	541	558	rBV	3236622	3055159	81.01%	10.044%
3	6.481	709	713	726	rBV	3073046	3072678	81.48%	10.101%
4	6.822	768	771	775	rBV	1046773	820429	21.76%	2.697%
5	7.392	863	868	871	rBV	2307223	2038834	54.06%	6.703%
6	8.104	985	989	993	rBV	1302335	1083978	28.74%	3.564%
7	9.180	1167	1172	1175	rBV	4755582	3771130	100.00%	12.398%
8	9.863	1284	1288	1291	rBV	1283973	1087028	28.82%	3.574%
9	10.410	1378	1381	1386	rVB	55297	51945	1.38%	0.171%
10	10.657	1419	1423	1426	rBV	2011821	1650091	43.76%	5.425%
11	10.757	1436	1440	1444	rBV4	34260	61830	1.64%	0.203%
12	11.351	1537	1541	1543	rBV	1011573	838467	22.23%	2.756%
13	11.374	1543	1545	1550	rVB	443755	350694	9.30%	1.153%
14	11.427	1551	1554	1557	rVV	118083	92361	2.45%	0.304%
15	11.592	1577	1582	1585	rBV3	42908	68349	1.81%	0.225%
16	11.727	1602	1605	1608	rBV	203302	147356	3.91%	0.484%
17	11.868	1623	1629	1636	rBV2	193979	384079	10.18%	1.263%
18	11.968	1642	1646	1648	rBV2	150686	163535	4.34%	0.538%
19	11.986	1648	1649	1654	rVB	66278	51780	1.37%	0.170%
20	12.145	1673	1676	1679	rBV	64109	58327	1.55%	0.192%
21	12.186	1680	1683	1686	rVB	50477	48570	1.29%	0.160%
22	12.415	1717	1722	1724	rBV3	297079	357457	9.48%	1.175%
23	12.433	1724	1725	1729	rVV3	258160	236747	6.28%	0.778%
24	12.498	1732	1736	1738	rBV3	71843	92311	2.45%	0.303%
25	12.521	1738	1740	1744	rVB2	90785	77011	2.04%	0.253%
26	12.568	1744	1748	1752	rBV	985321	862470	22.87%	2.835%
27	12.657	1760	1763	1768	rVB3	65592	67772	1.80%	0.223%
28	12.798	1781	1787	1793	rBV	891956	1001712	26.56%	3.293%
29	12.851	1793	1796	1799	rVB	68423	68283	1.81%	0.224%
30	12.939	1806	1811	1814	rBV	2763628	2433299	64.52%	7.999%
31	13.021	1820	1825	1829	rBV3	80217	140071	3.71%	0.460%
32	13.127	1836	1843	1846	rBV	175306	267594	7.10%	0.880%
33	13.192	1851	1854	1857	rBV	116929	101846	2.70%	0.335%
34	13.233	1858	1861	1865	rVV2	117358	108834	2.89%	0.358%
35	13.327	1874	1877	1879	rBV	98445	83244	2.21%	0.274%
36	13.357	1879	1882	1891	rVB2	97686	135090	3.58%	0.444%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.645	1928	1931	1934	rBV	96987	88449	2.35%	0.291%
38	13.751	1947	1949	1951	rBV	83746	61433	1.63%	0.202%
39	13.774	1951	1953	1956	rVB	90964	75423	2.00%	0.248%
40	13.804	1956	1958	1960	rBV	73181	77778	2.06%	0.256%
41	14.004	1987	1992	1994	rBV2	1083565	1371747	36.37%	4.510%
42	14.027	1994	1996	2003	rVB	522595	507114	13.45%	1.667%
43	14.109	2008	2010	2012	rVB	91620	63556	1.69%	0.209%
44	14.392	2053	2058	2060	rBV2	212901	268736	7.13%	0.883%
45	14.427	2061	2064	2065	rBV	104520	101749	2.70%	0.334%
46	14.892	2140	2143	2147	rBV	288073	293169	7.77%	0.964%
47	15.051	2166	2170	2173	rBV	496447	720210	19.10%	2.368%
48	15.356	2219	2222	2226	rVB	281523	311627	8.26%	1.024%
49	15.415	2229	2232	2236	rVB	398413	460388	12.21%	1.514%
50	15.480	2240	2243	2246	rVB	488635	511379	13.56%	1.681%
51	16.962	2492	2495	2503	rVB3	195062	369657	9.80%	1.215%

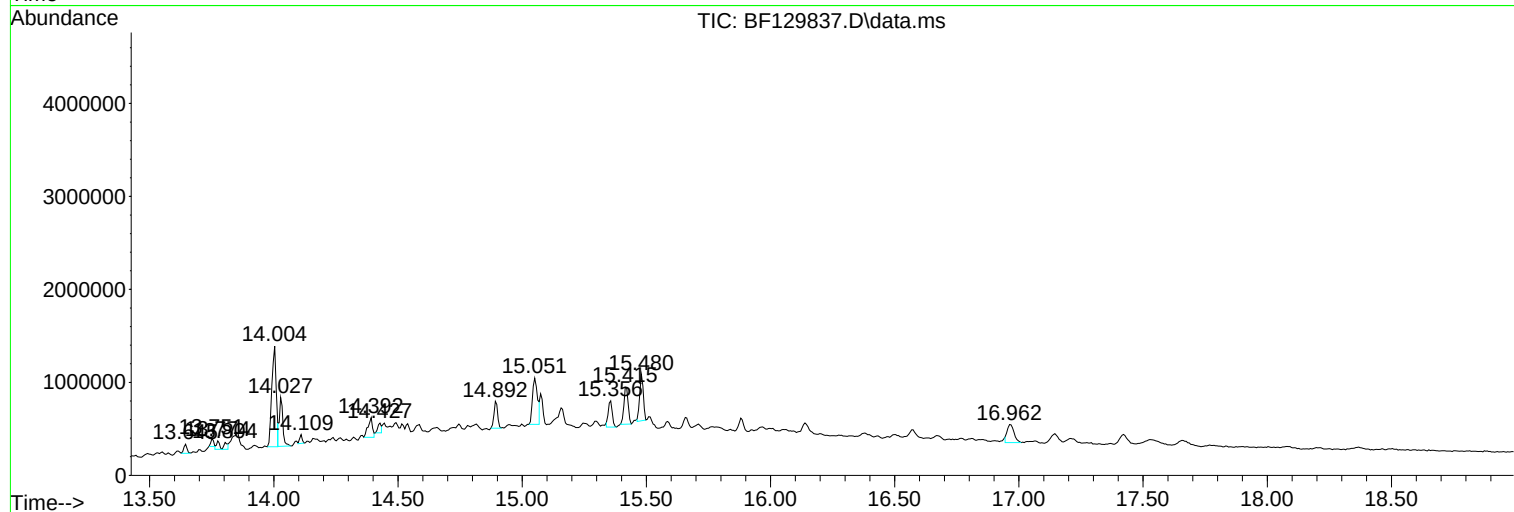
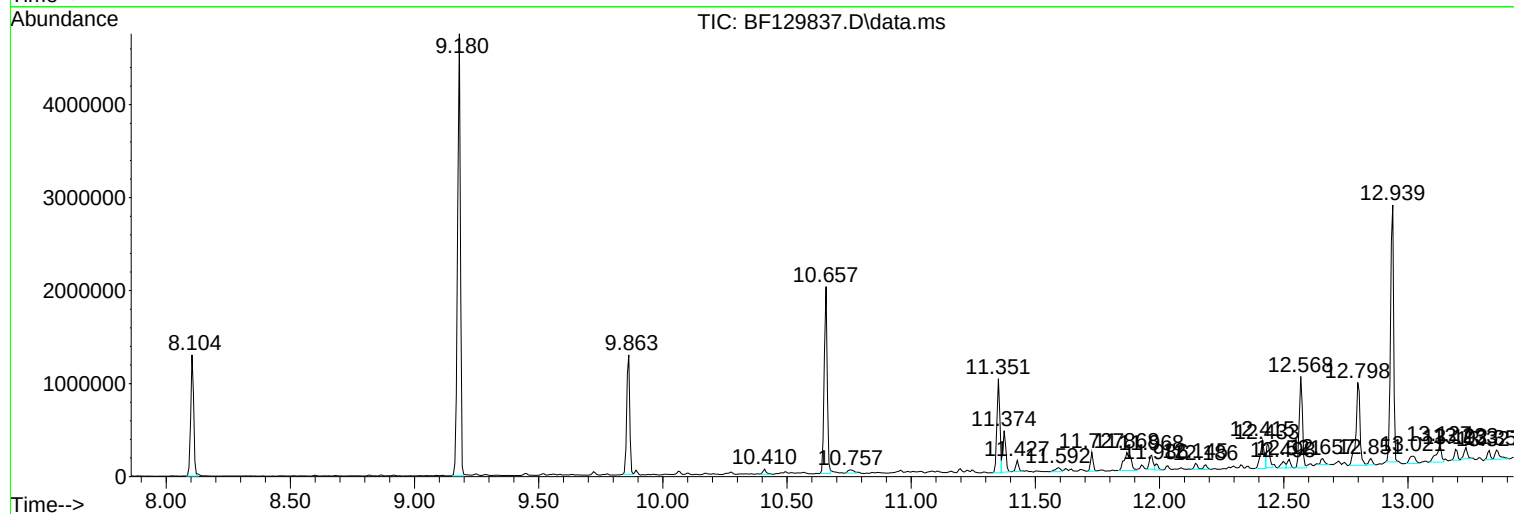
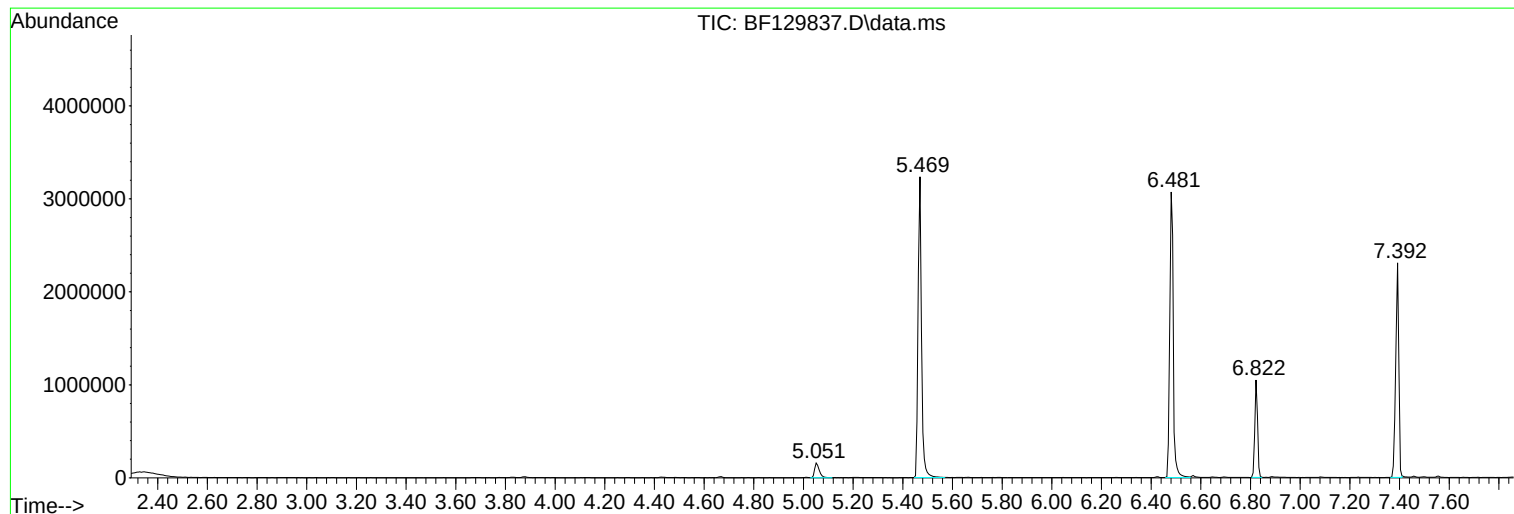
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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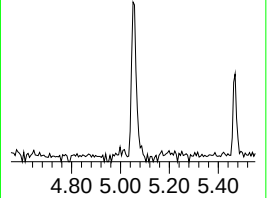
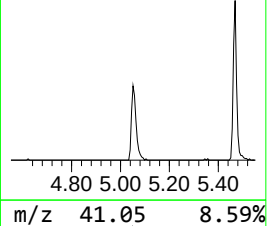
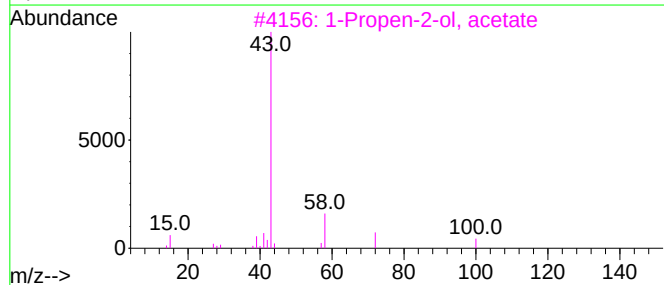
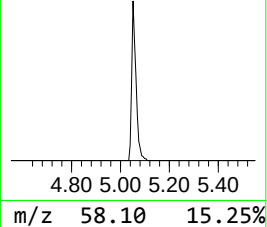
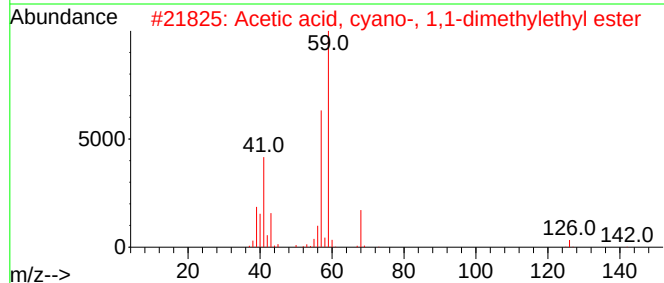
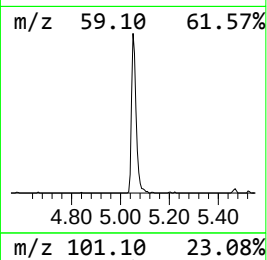
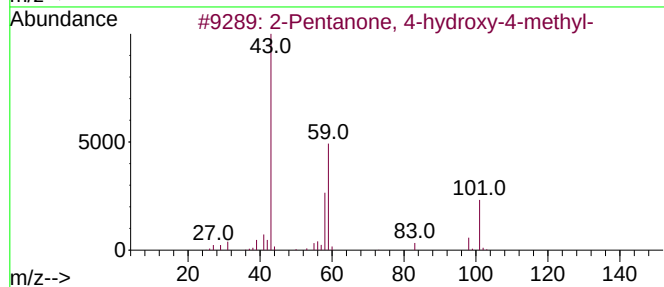
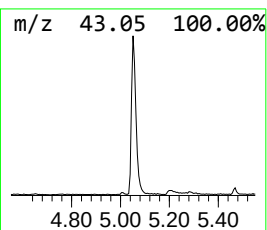
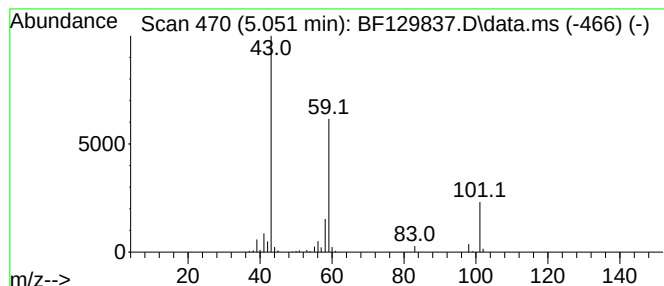
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.051	5.01 ng	205652	1,4-Dichlorobenzene-d4	6.822

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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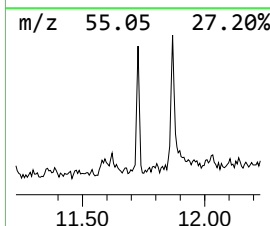
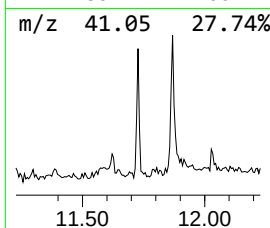
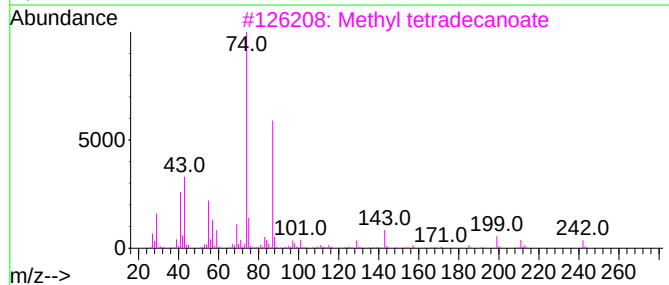
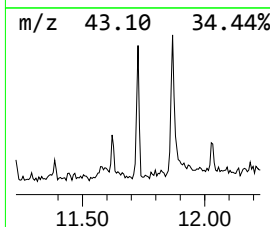
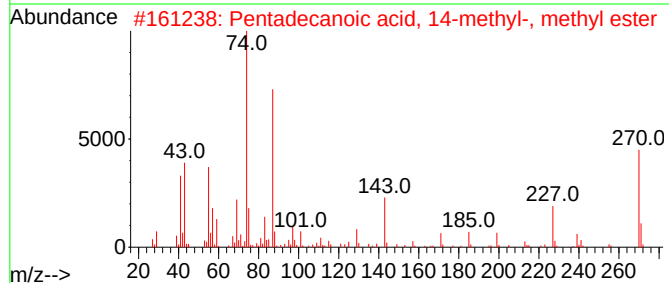
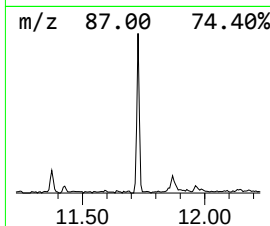
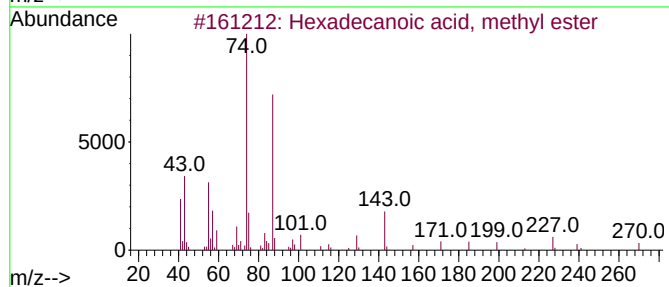
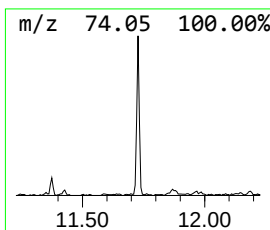
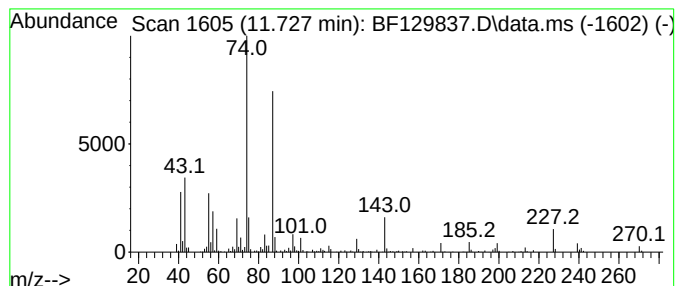
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexadecanoic acid, methyl e... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.727	3.51 ng	147356	Phenanthrene-d10	11.351

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	94
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5			Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	83



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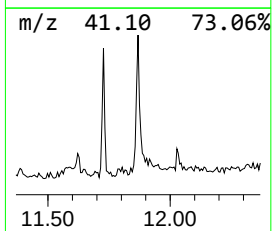
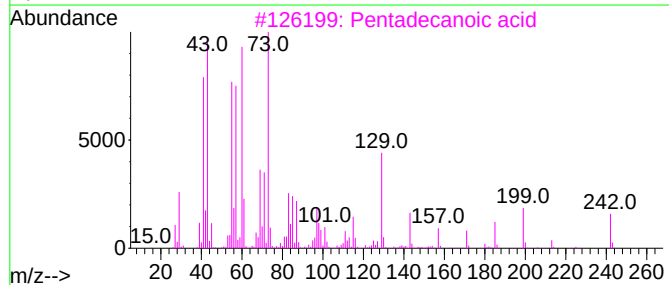
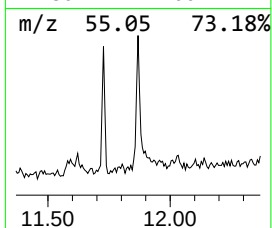
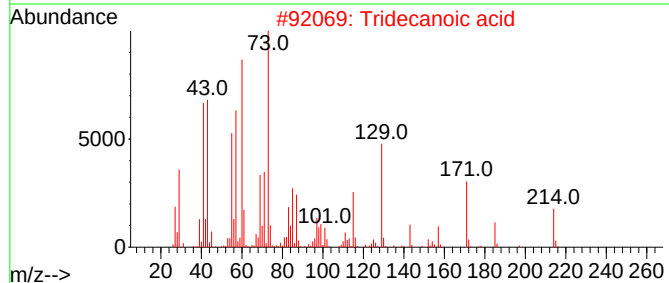
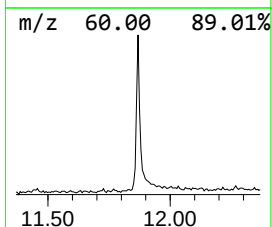
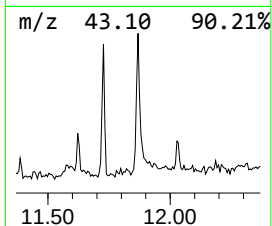
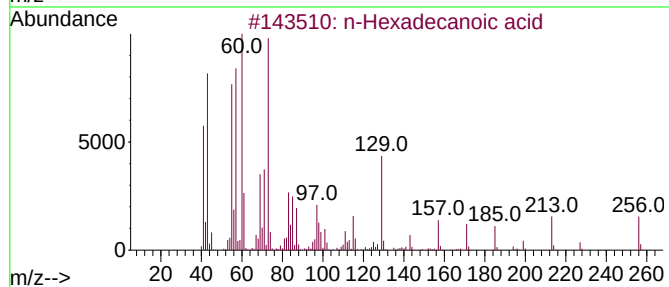
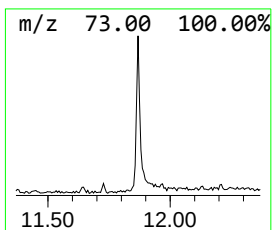
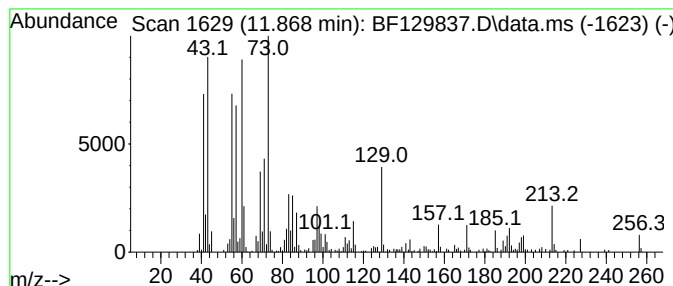
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.868	9.16 ng	384079	Phenanthrene-d10		11.351	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	98
2	Tridecanoic acid		214	C13H26O2	000638-53-9	92
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	91
4	n-Decanoic acid		172	C10H20O2	000334-48-5	76
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	58



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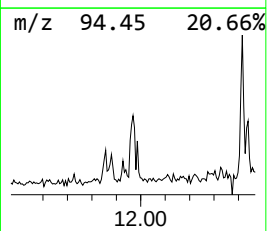
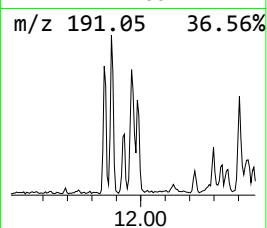
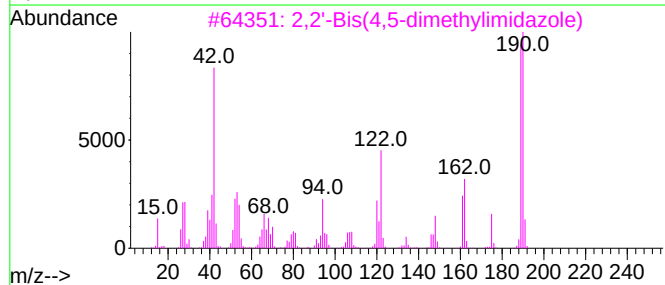
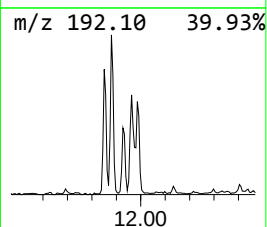
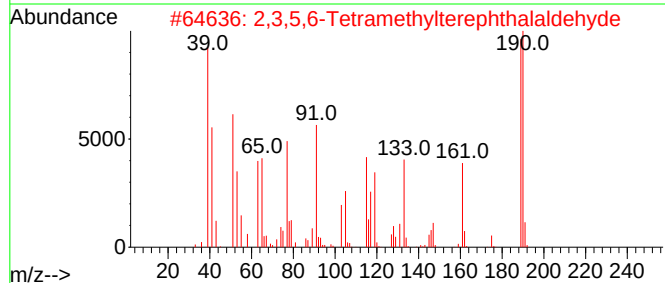
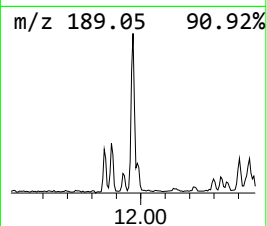
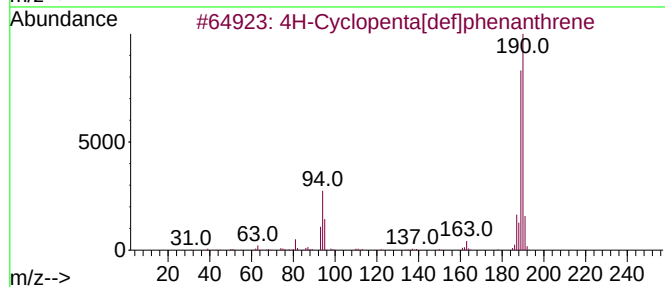
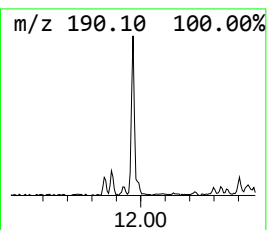
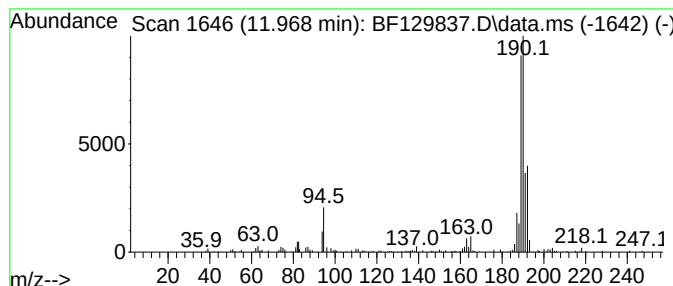
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.968	3.90 ng	163535	Phenanthrene-d10	11.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
4			Methyl diselenide	190	C2H6Se2	007101-31-7	38
5			Pyrimidine, 6-oxo-4-(3-fluorophe...	190	C10H7FN2O	085979-55-1	38



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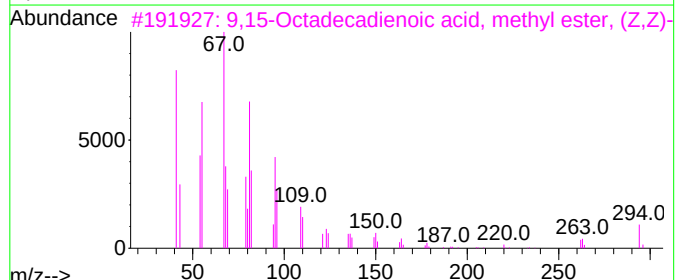
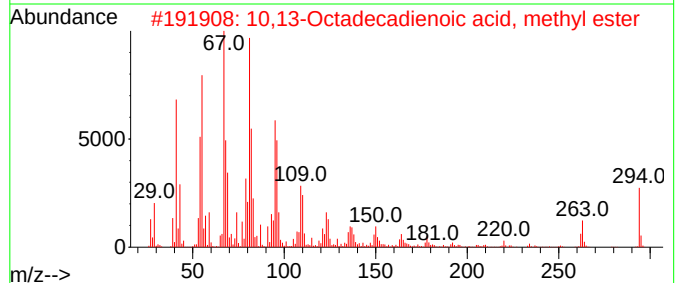
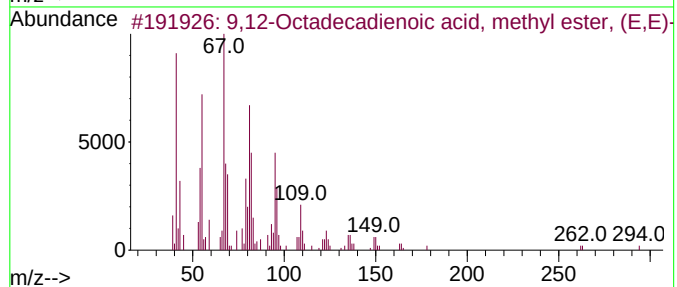
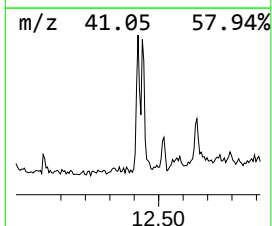
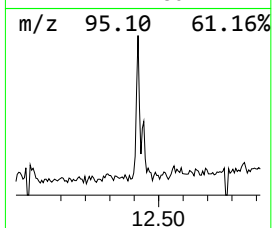
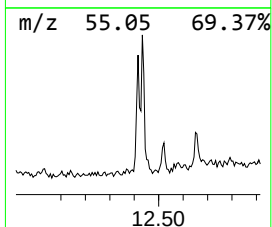
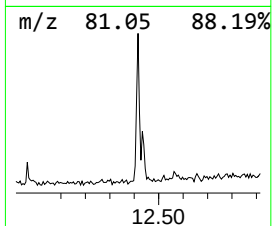
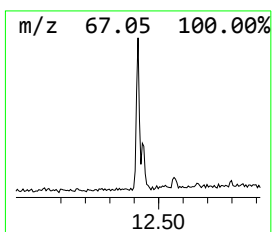
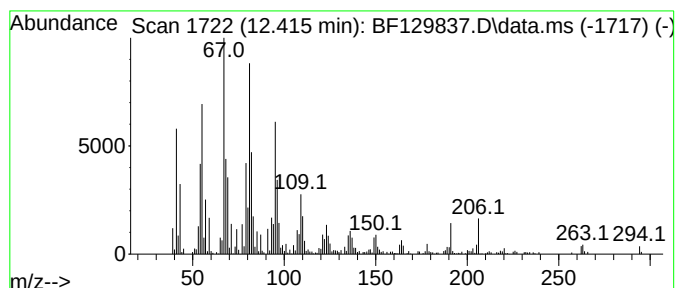
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ClientSampleId :
HD-02-081222

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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 9,12-Octadecadienoic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.415	8.53 ng	357457	Phenanthrene-d10		11.351	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...		294	C19H34O2	002566-97-4	99
2	10,13-Octadecadienoic acid, meth...		294	C19H34O2	056554-62-2	99
3	9,15-Octadecadienoic acid, methy...		294	C19H34O2	017309-05-6	99
4	9,12-Octadecadienoic acid (Z,Z)-...		294	C19H34O2	000112-63-0	99
5	9,12-Octadecadienoic acid, methy...		294	C19H34O2	002462-85-3	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

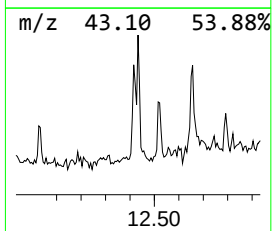
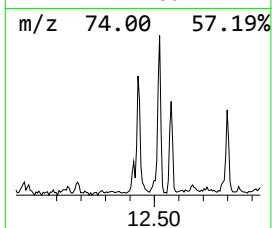
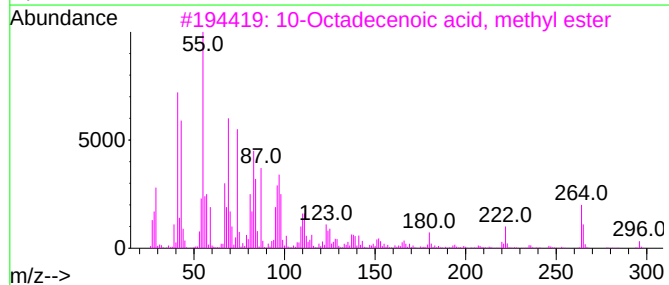
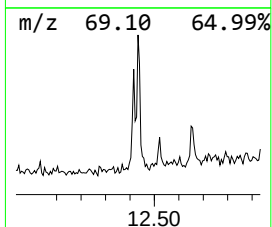
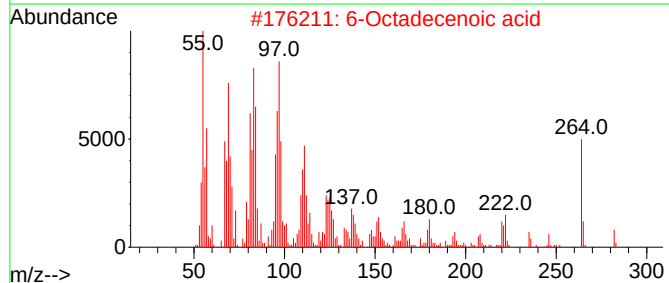
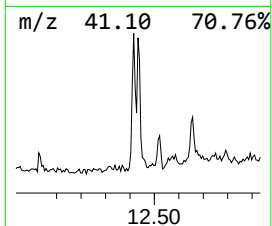
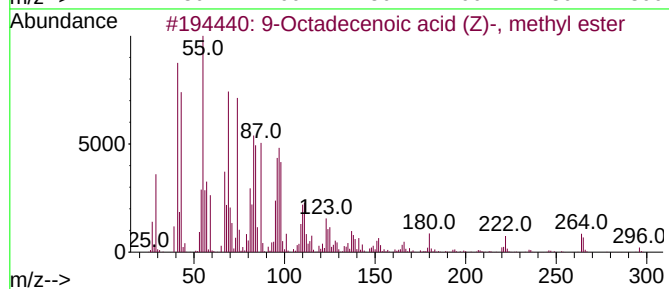
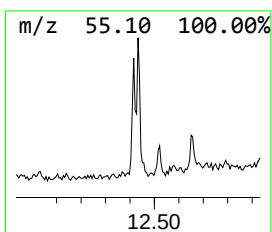
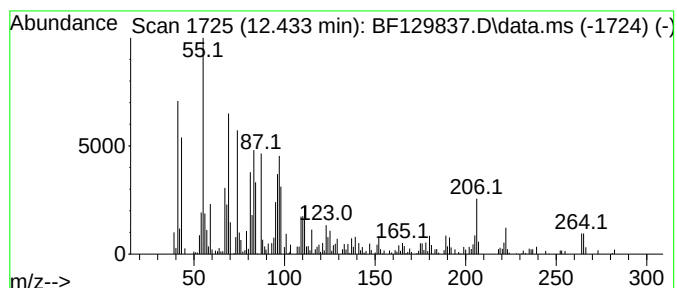
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9-Octadecenoic acid (Z)-, m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.433	5.65 ng	236747	Phenanthrene-d10	11.351	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	93
2	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	91
3	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	90
4	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	90
5	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
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Misc :
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Instrument :
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ClientSampleId :
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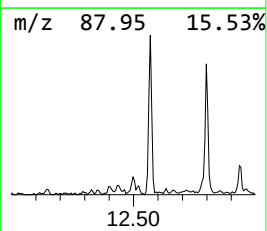
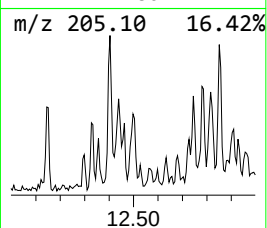
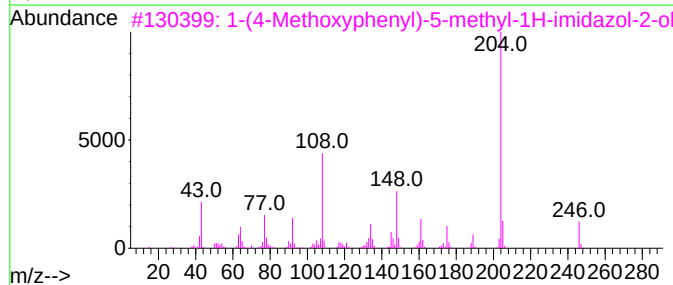
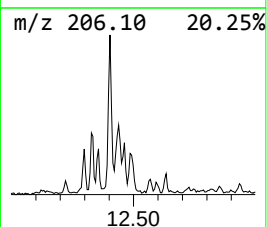
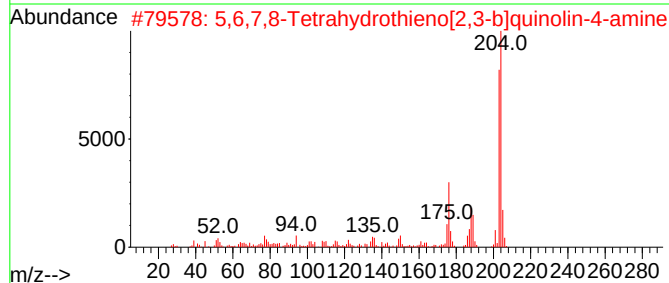
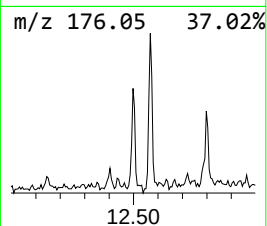
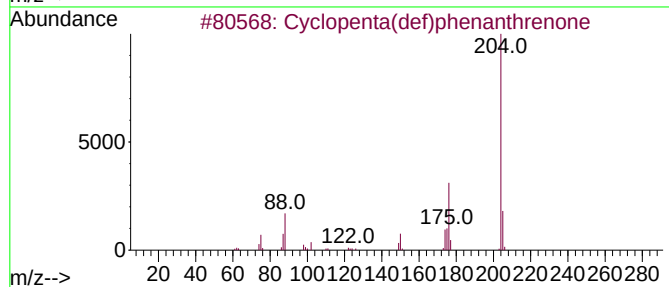
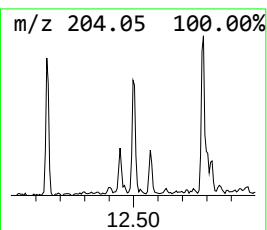
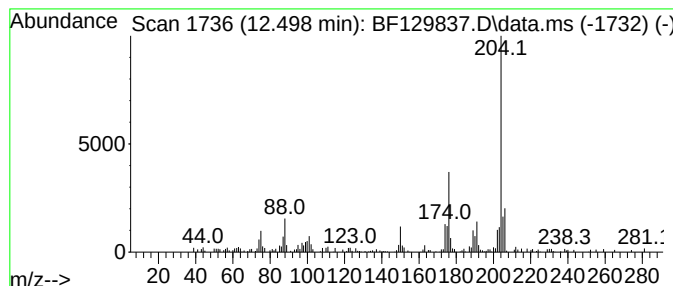
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.498	2.20 ng	92311	Phenanthrene-d10	11.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	58
3			1-(4-Methoxyphenyl)-5-methyl-1H-...	246	C13H14N2O3	1010500-96-8	47
4			6-Chloro-2,3,4,9-tetrahydro-1H-c...	248	C13H13ClN2O	1000303-44-7	47
5			1-(4-Methoxyphenyl)-5-methyl-2,3...	204	C11H12N2O2	014058-90-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

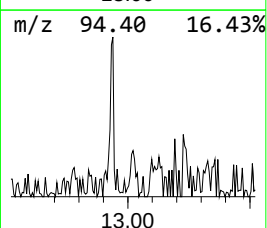
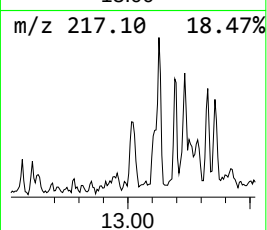
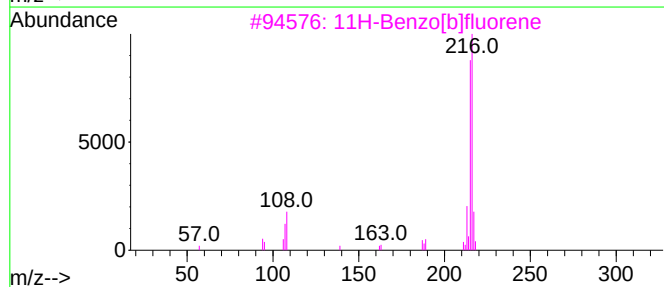
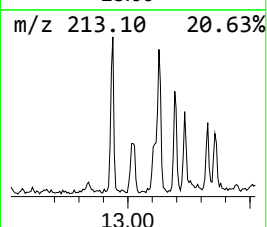
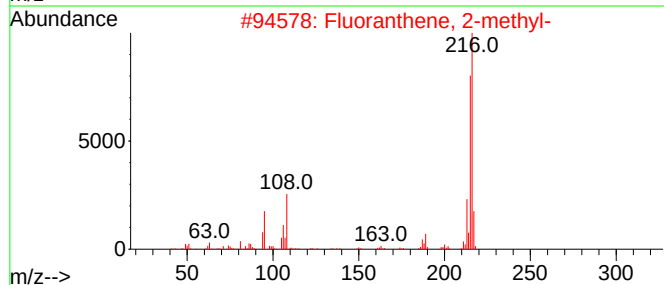
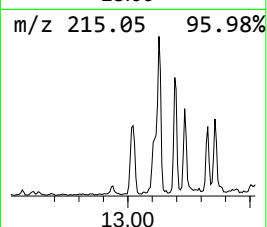
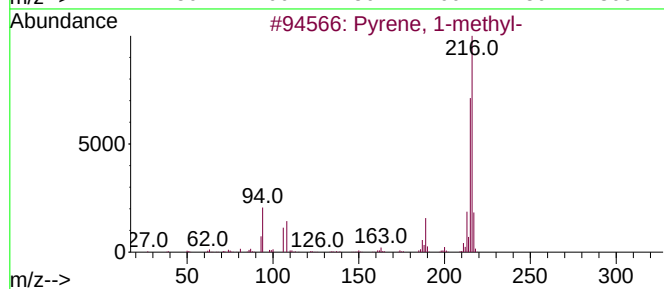
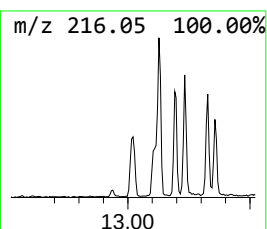
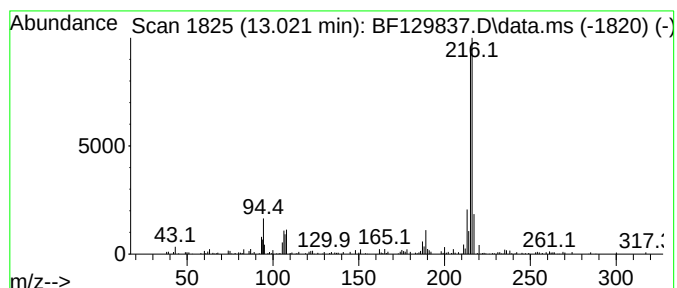
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.021	2.04 ng	140071	Chrysene-d12	14.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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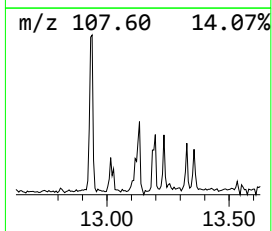
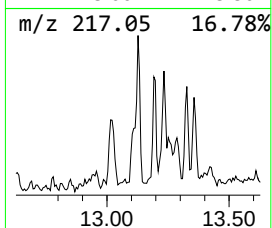
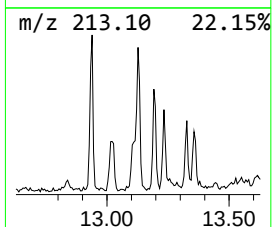
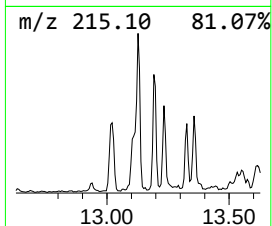
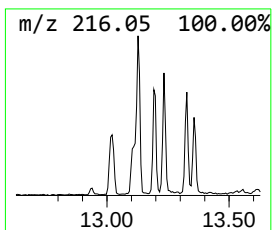
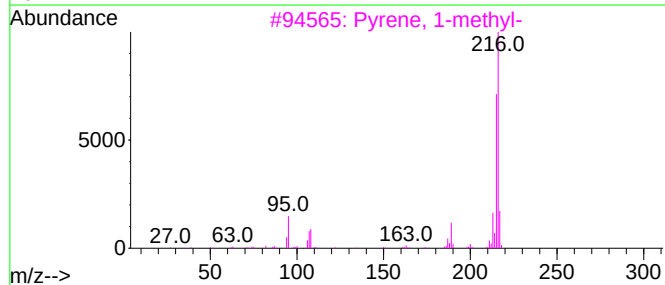
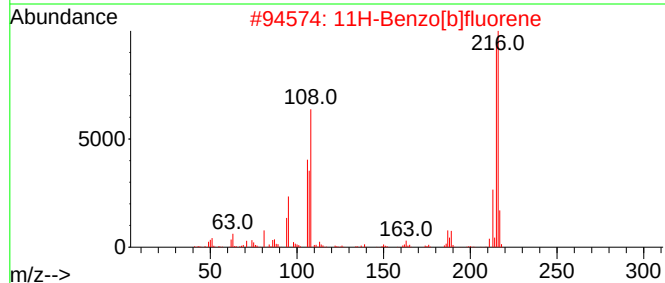
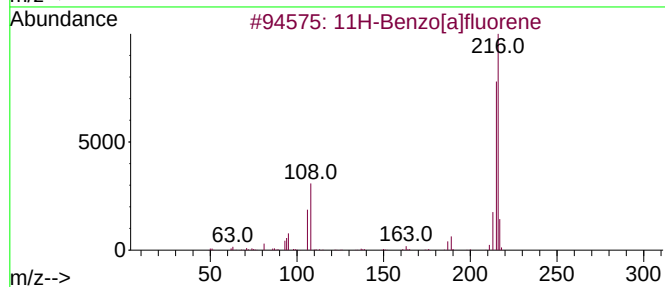
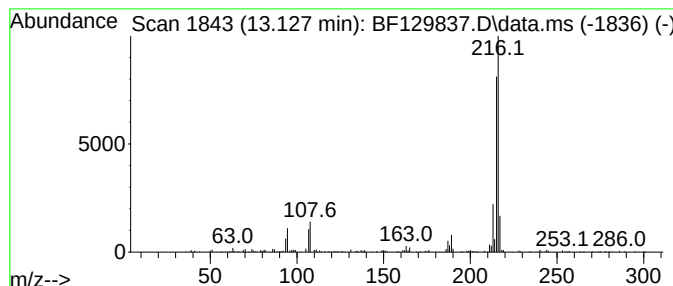
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.127	3.90 ng	267594	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-02-081222

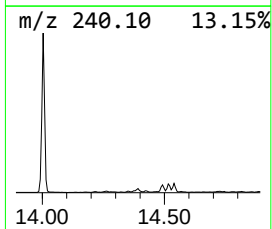
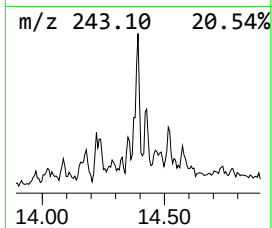
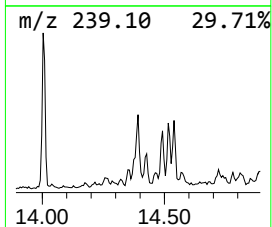
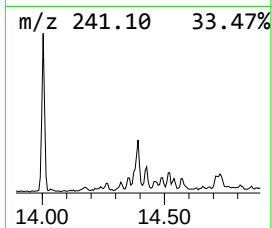
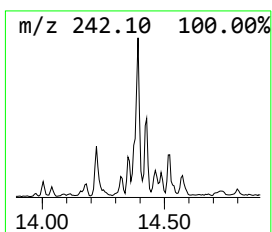
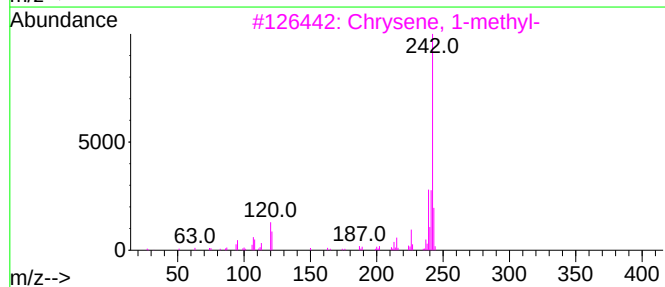
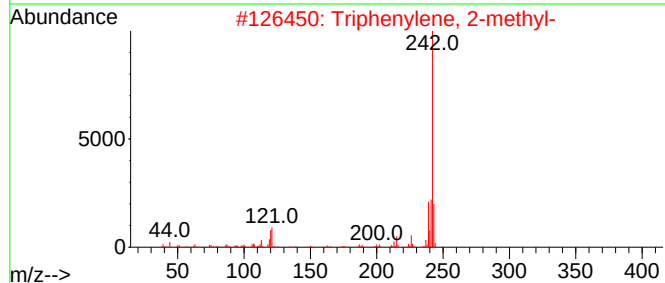
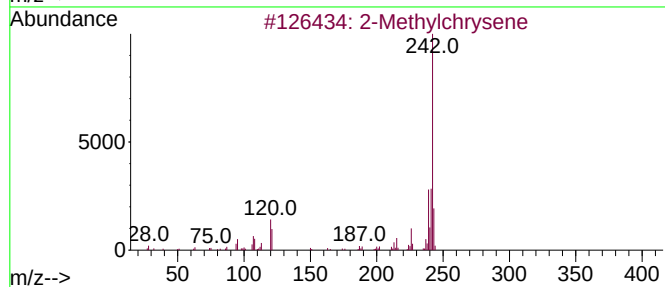
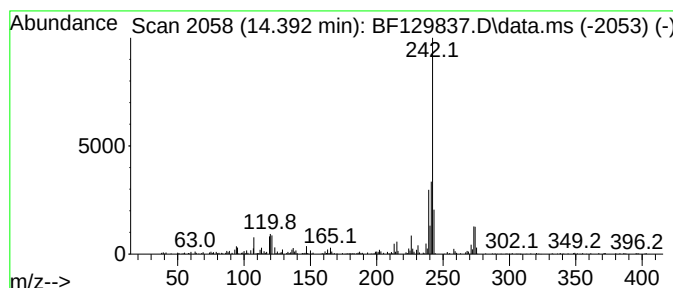
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Methylchrysene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.392	3.92 ng	268736	Chrysene-d12	14.004

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylchrysene	242	C19H14	003351-32-4	96	
2	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	95	
3	Chrysene, 1-methyl-	242	C19H14	003351-28-8	95	
4	Benz[a]anthracene, 12-methyl-	242	C19H14	002422-79-9	91	
5	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	91	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
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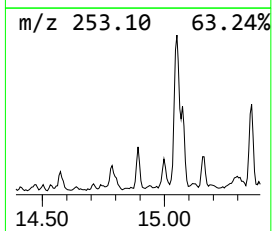
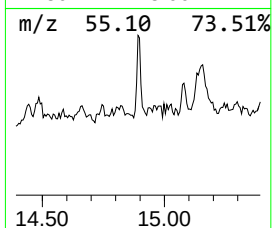
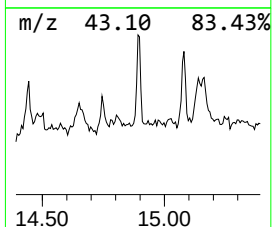
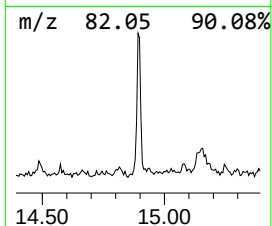
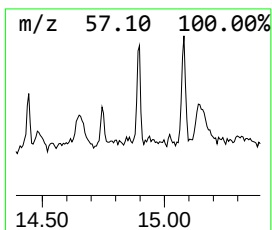
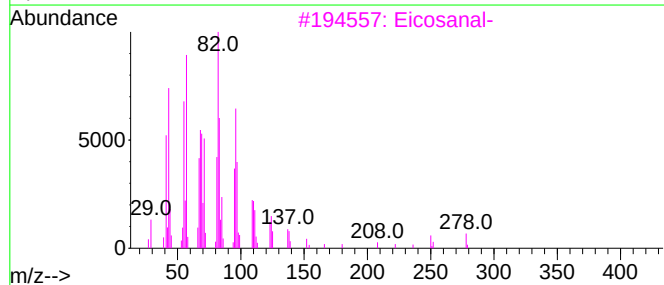
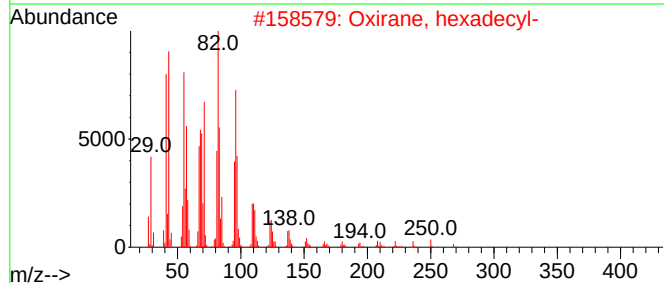
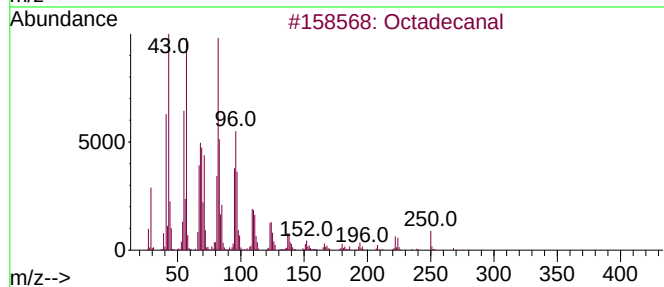
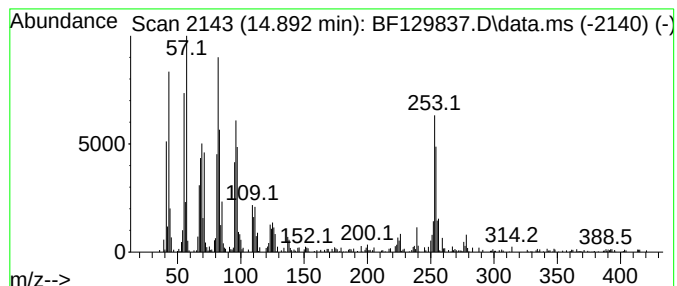
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.892	11.47 ng	293169	Perylene-d12	15.480	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal	268	C18H36O	000638-66-4	92
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3	Eicosanal-	296	C20H40O	002400-66-0	90
4	17-Octadecenal	266	C18H34O	056554-86-0	89
5	4,5-Dihydrobenzo[e]pyrene	254	C20H14	095676-42-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-081222

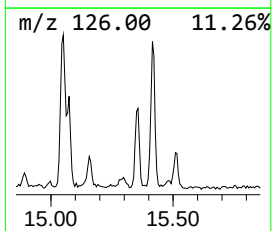
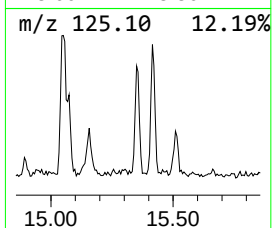
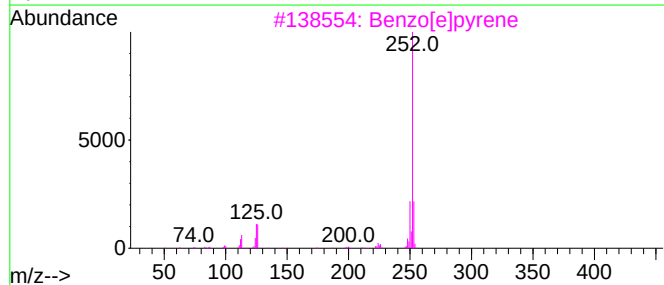
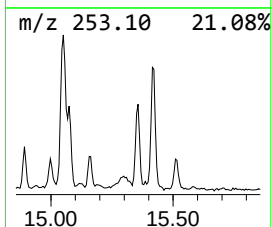
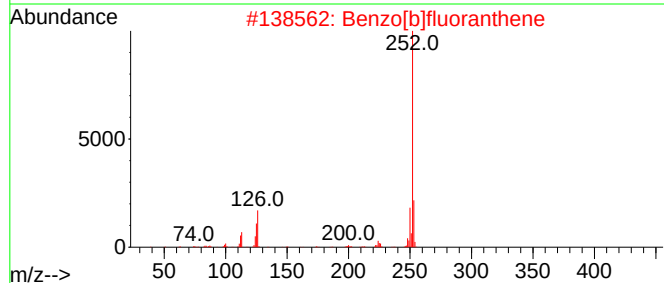
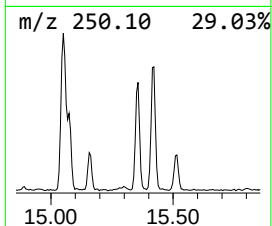
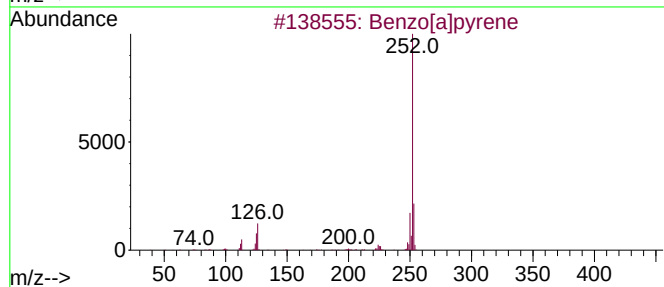
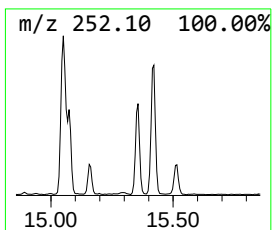
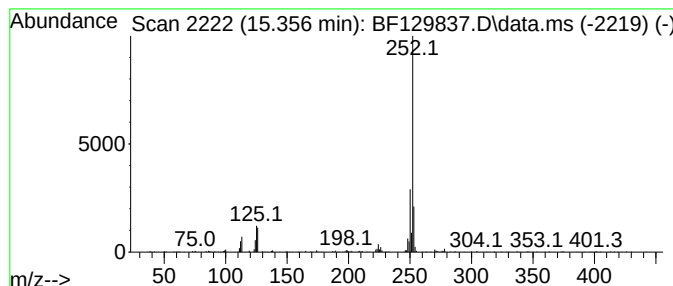
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.356	12.19 ng	311627	Perylene-d12	15.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	97
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
3			Benzo[e]pyrene	252	C20H12	000192-97-2	94
4			Perylene	252	C20H12	000198-55-0	90
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
Data File : BF129837.D
Acq On : 17 Aug 2022 10:51
Operator : CG\JU
Sample : N4199-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-081222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.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
2-Pentanone, 4-...	5.051	5.0	ng	205652	1	6.822	820429	20.0
Hexadecanoic ac...	11.727	3.5	ng	147356	4	11.351	838467	20.0
n-Hexadecanoic ...	11.868	9.2	ng	384079	4	11.351	838467	20.0
4H-Cyclopenta[d...	11.968	3.9	ng	163535	4	11.351	838467	20.0
9,12-Octadecadi...	12.415	8.5	ng	357457	4	11.351	838467	20.0
9-Octadecenoic ...	12.433	5.7	ng	236747	4	11.351	838467	20.0
Cyclopenta(def)...	12.498	2.2	ng	92311	4	11.351	838467	20.0
Pyrene, 1-methyl-	13.021	2.0	ng	140071	5	14.004	1371750	20.0
11H-Benzo[a]flu...	13.127	3.9	ng	267594	5	14.004	1371750	20.0
2-Methylchrysene	14.392	3.9	ng	268736	5	14.004	1371750	20.0
Octadecanal	14.892	11.5	ng	293169	6	15.480	511379	20.0
Benzo[e]pyrene	15.356	12.2	ng	311627	6	15.480	511379	20.0