

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
 Data File : BF127267.D
 Acq On : 08 Mar 2022 19:28
 Operator : CG\JU
 Sample : N1829-01 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-030722

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-BF021622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127267.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.493	542	545	554	rBV	455241	460147	12.68%	2.293%
2	6.510	714	718	726	rBV	464063	468865	12.92%	2.336%
3	6.881	777	781	784	rBV	775193	695052	19.16%	3.464%
4	7.440	872	876	879	rBV	377948	299725	8.26%	1.494%
5	8.163	995	999	1002	rBV	1079142	859529	23.69%	4.283%
6	9.234	1178	1181	1188	rVB	626183	482012	13.28%	2.402%
7	9.922	1294	1298	1301	rBV	862856	669790	18.46%	3.338%
8	9.951	1301	1303	1308	rVB	74931	61667	1.70%	0.307%
9	10.339	1361	1369	1372	rBV4	27642	38437	1.06%	0.192%
10	10.469	1385	1391	1397	rVB	54648	54298	1.50%	0.271%
11	10.710	1429	1432	1436	rBV	227805	206563	5.69%	1.029%
12	10.810	1447	1449	1453	rBV3	23321	38436	1.06%	0.192%
13	11.416	1548	1552	1554	rBV	670978	623211	17.18%	3.106%
14	11.439	1554	1556	1559	rVV	432146	338026	9.32%	1.684%
15	11.486	1562	1564	1567	rVB	108081	87309	2.41%	0.435%
16	11.645	1584	1591	1595	rBV	50639	61723	1.70%	0.308%
17	11.792	1613	1616	1620	rBV	1156477	900597	24.82%	4.488%
18	11.916	1632	1637	1639	rBV	64806	62302	1.72%	0.310%
19	11.945	1639	1642	1645	rVB	81576	74453	2.05%	0.371%
20	12.027	1653	1656	1659	rBV2	181760	182910	5.04%	0.911%
21	12.210	1681	1687	1690	rBV2	55964	63429	1.75%	0.316%
22	12.239	1690	1692	1697	rVB2	55665	45853	1.26%	0.228%
23	12.486	1726	1734	1736	rBV	3744808	3628256	100.00%	18.080%
24	12.504	1736	1737	1743	rVV2	2210721	1830230	50.44%	9.120%
25	12.557	1743	1746	1748	rVV	70504	72663	2.00%	0.362%
26	12.586	1748	1751	1755	rVB	487184	383206	10.56%	1.910%
27	12.633	1755	1759	1762	rBV	1398117	1142386	31.49%	5.693%
28	12.786	1777	1785	1787	rBV2	52384	91755	2.53%	0.457%
29	12.863	1791	1798	1803	rVV	1108024	1222461	33.69%	6.092%
30	12.910	1803	1806	1810	rVB2	46070	50338	1.39%	0.251%
31	12.998	1815	1821	1823	rBV2	404099	360520	9.94%	1.797%
32	13.022	1823	1825	1828	rVB	48633	45640	1.26%	0.227%
33	13.074	1830	1834	1838	rBV2	60843	108130	2.98%	0.539%
34	13.186	1846	1853	1856	rBV	163276	231685	6.39%	1.155%
35	13.257	1862	1865	1867	rVB	129042	103056	2.84%	0.514%
36	13.292	1867	1871	1873	rVV2	76576	90642	2.50%	0.452%

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37	13.316	1873	1875	1879	rVB2	71047	61089	1.68%	0.304%
38	13.386	1884	1887	1890	rBV	55091	55722	1.54%	0.278%
39	13.421	1890	1893	1899	rVB3	69326	95157	2.62%	0.474%
40	13.698	1938	1940	1944	rVB2	31845	37364	1.03%	0.186%
41	13.810	1956	1959	1961	rBV	73815	66102	1.82%	0.329%
42	13.839	1961	1964	1966	rVV	65898	60512	1.67%	0.302%
43	13.863	1966	1968	1974	rVV2	73512	93328	2.57%	0.465%
44	13.980	1986	1988	1992	rVB2	54351	47992	1.32%	0.239%
45	14.063	1997	2002	2004	rVV2	640011	921763	25.41%	4.593%
46	14.086	2004	2006	2013	rVB	440955	409888	11.30%	2.043%
47	14.168	2014	2020	2023	rBV2	145649	197425	5.44%	0.984%
48	14.598	2091	2093	2096	rVB2	66525	56273	1.55%	0.280%
49	15.115	2177	2181	2184	rBV	309844	449045	12.38%	2.238%
50	15.427	2231	2234	2240	rVB	184183	231240	6.37%	1.152%
51	15.492	2241	2245	2251	rVB	312230	397244	10.95%	1.980%
52	15.557	2252	2256	2259	rBV	257666	331010	9.12%	1.650%
53	15.586	2259	2261	2268	rVB	83288	112513	3.10%	0.561%
54	17.074	2509	2514	2520	rVB2	95292	178460	4.92%	0.889%
55	17.533	2588	2592	2598	rBV	76437	129821	3.58%	0.647%

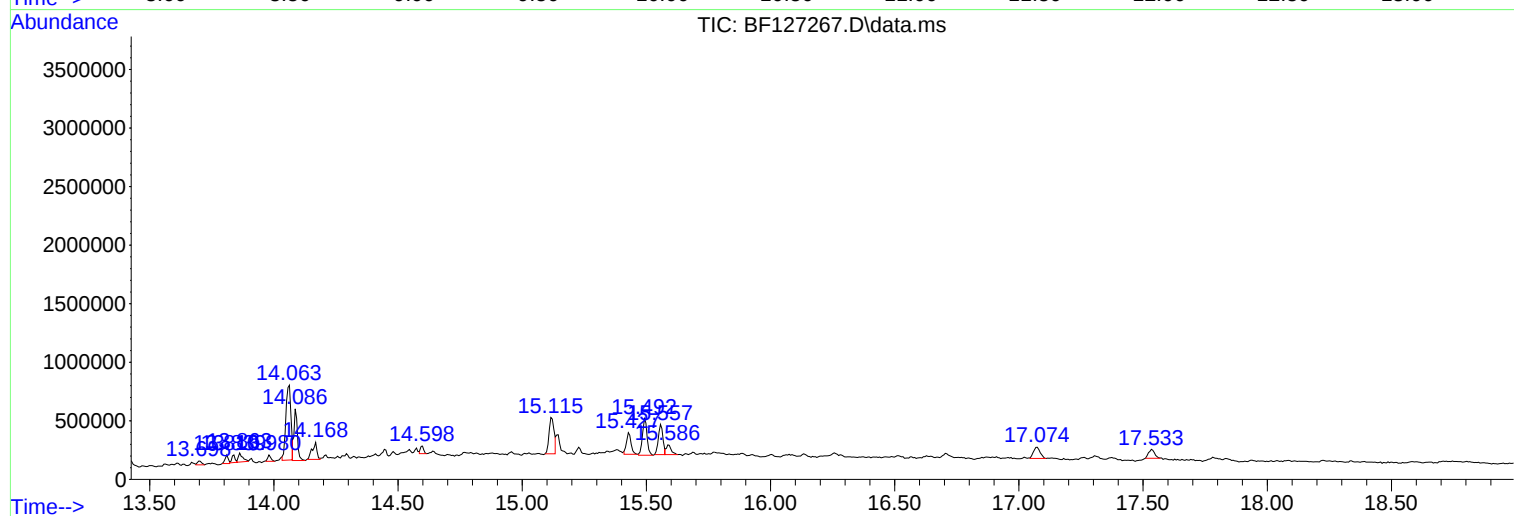
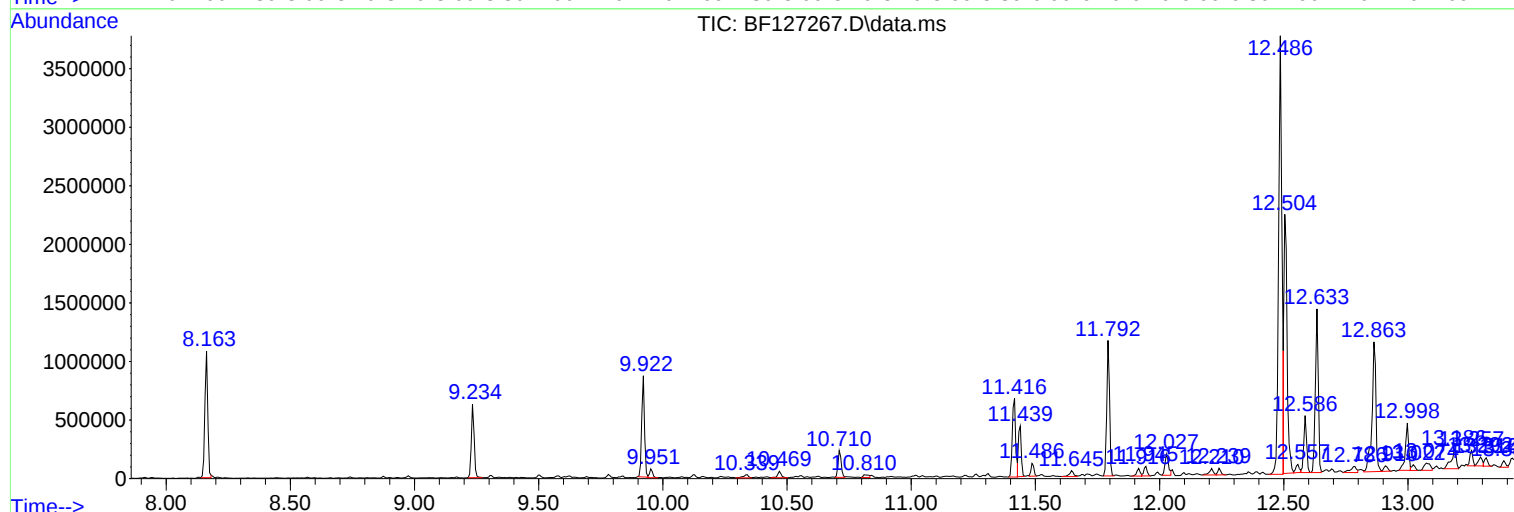
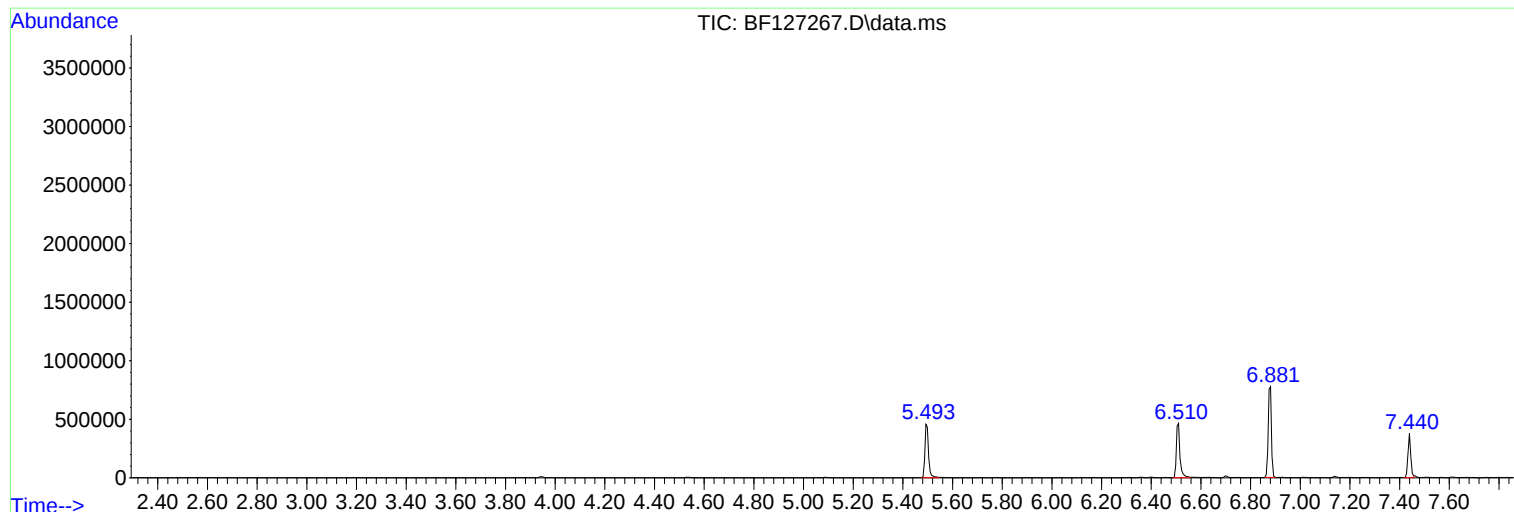
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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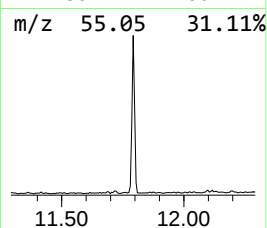
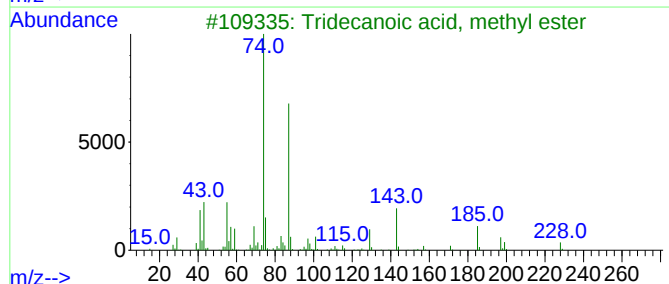
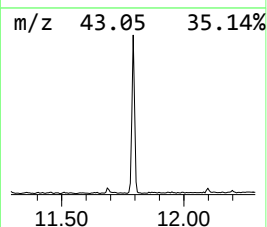
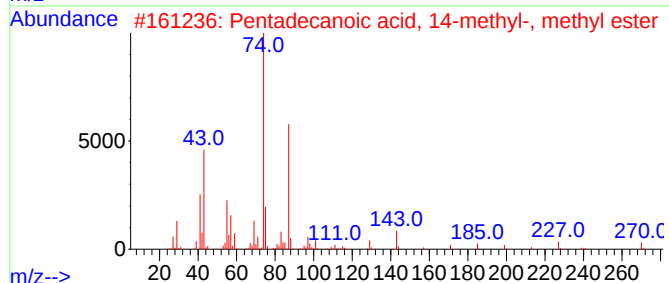
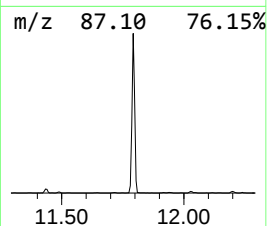
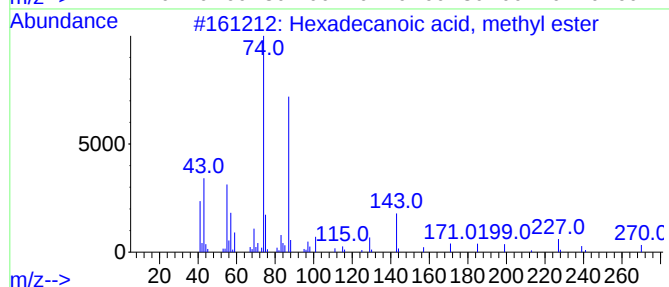
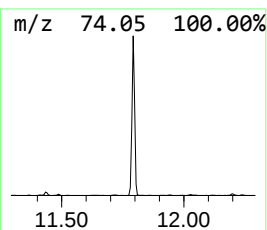
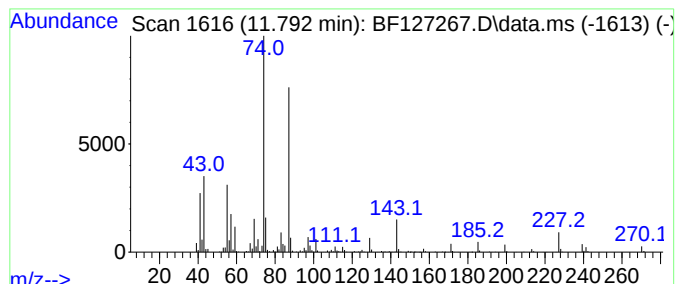
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.792	28.90 ng	900597	Phenanthrene-d10	11.416

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	95
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	83



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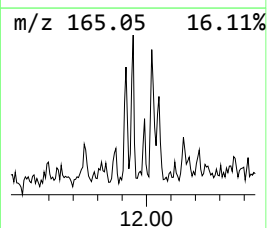
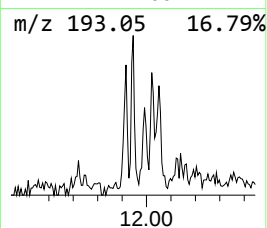
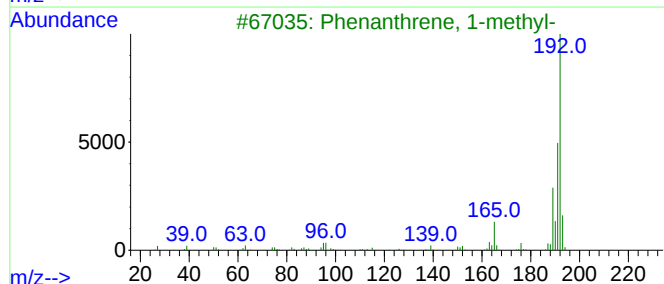
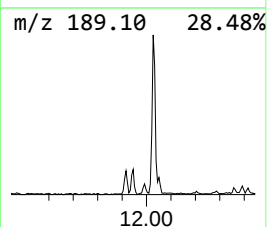
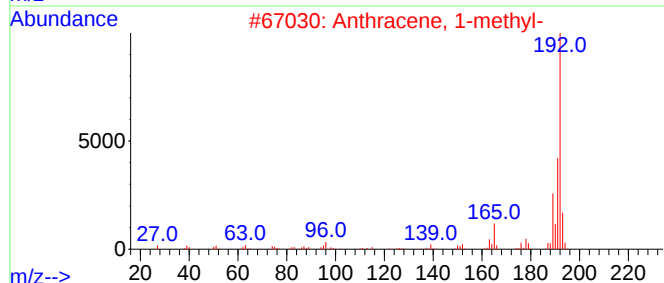
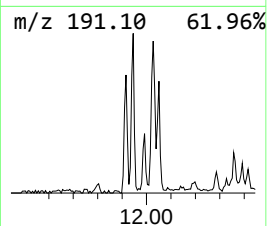
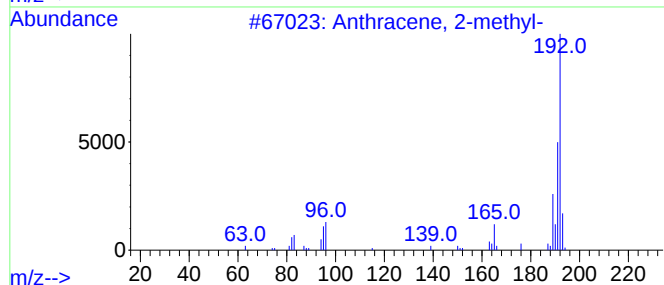
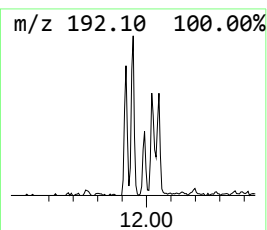
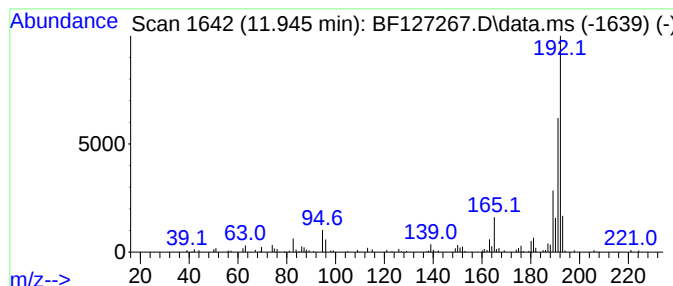
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	2.39 ng	74453	Phenanthrene-d10	11.416

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90



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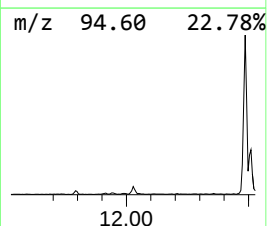
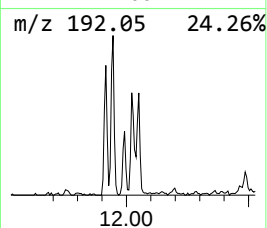
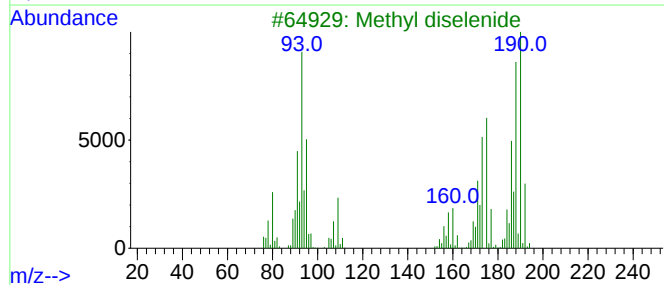
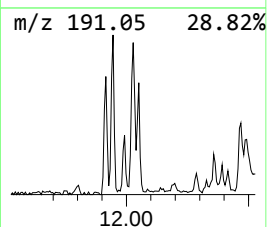
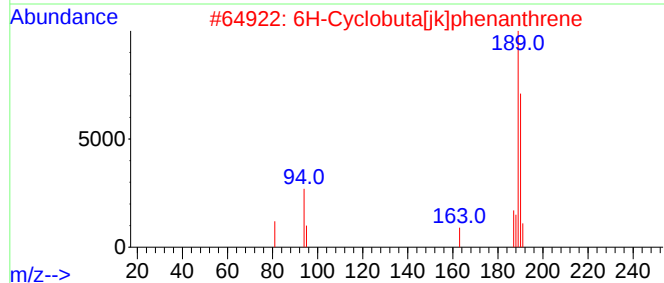
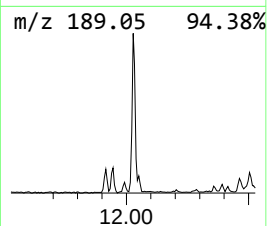
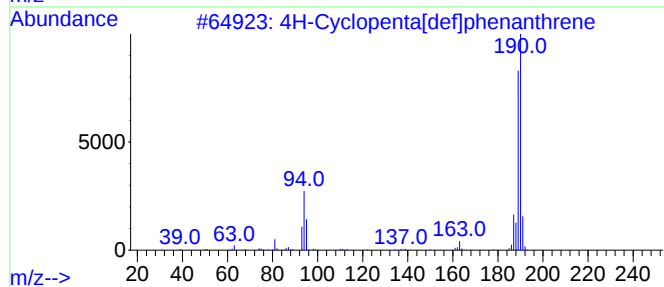
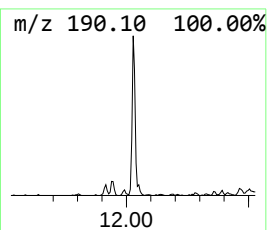
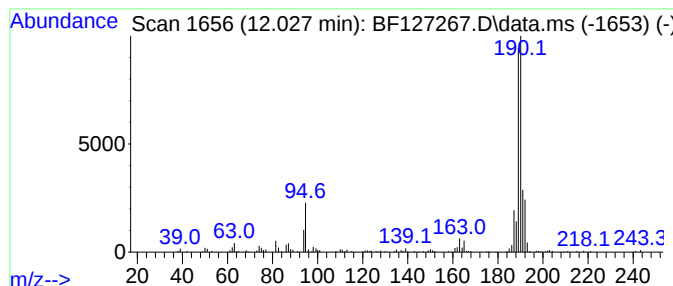
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.027	5.87 ng	182910	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	64
3	Methyl diselenide		190	C2H6Se2	007101-31-7	49
4	2,2'-Bis(4,5-dimethylimidazole)		190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...		190	C10H6O4	004923-55-1	25



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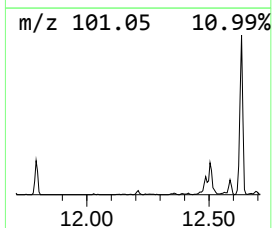
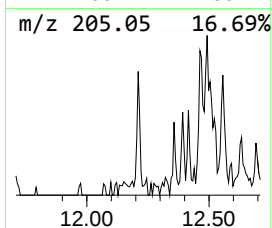
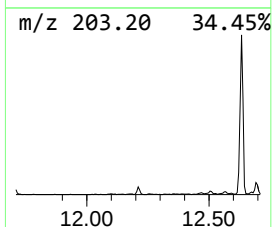
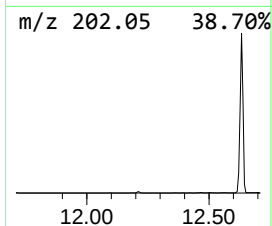
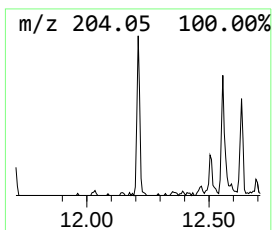
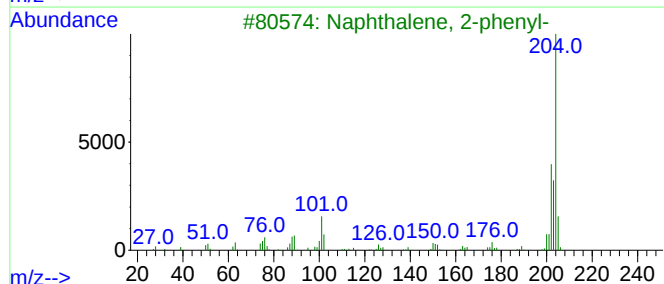
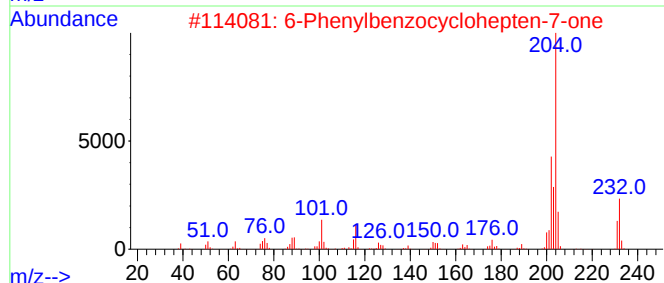
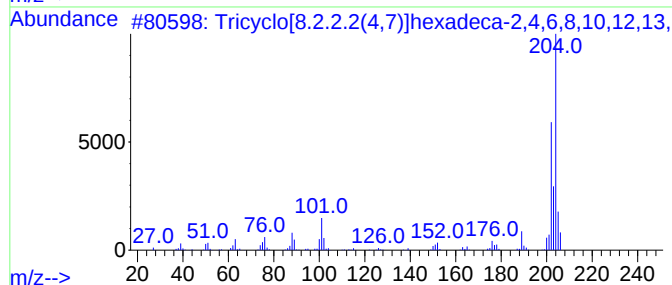
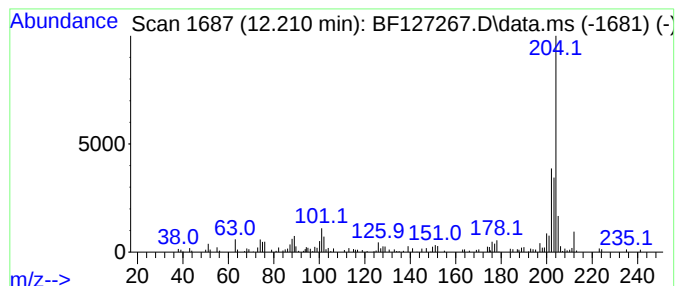
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.210	2.04 ng	63429	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	83
2			6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	80
3			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	78
4			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	55
5			9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	49



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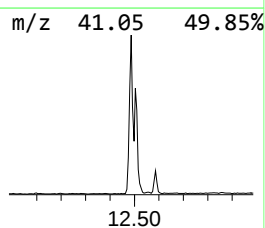
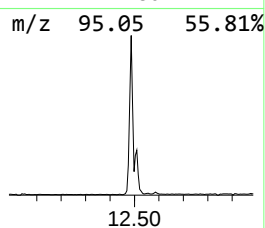
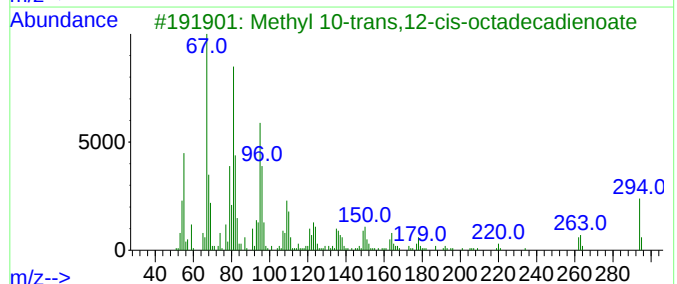
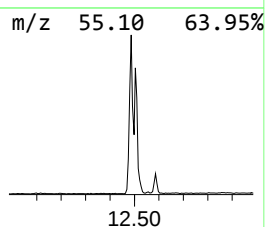
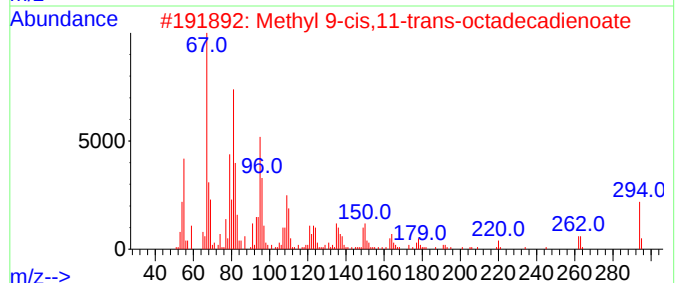
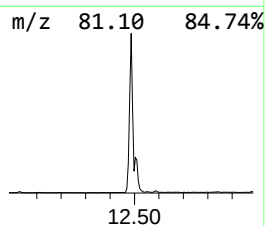
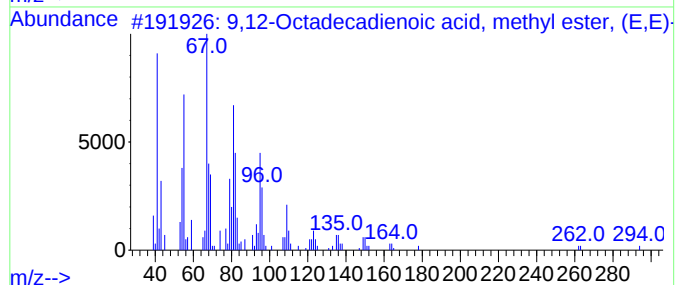
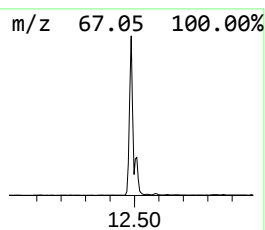
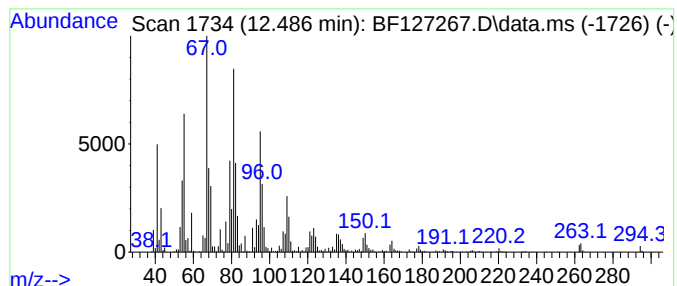
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	116.44 ng	3628260	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	99
3			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
4			9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5			9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-030722

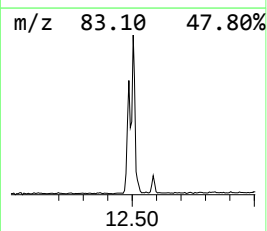
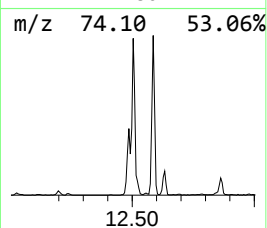
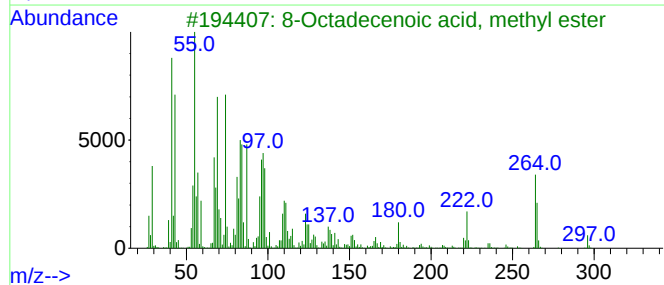
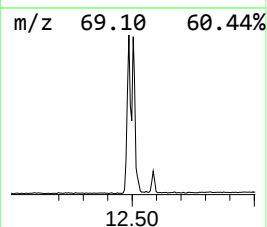
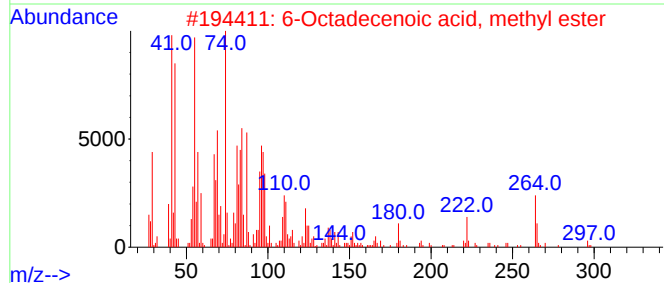
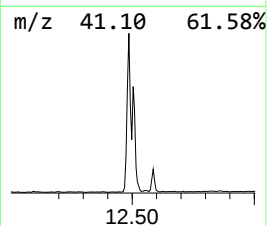
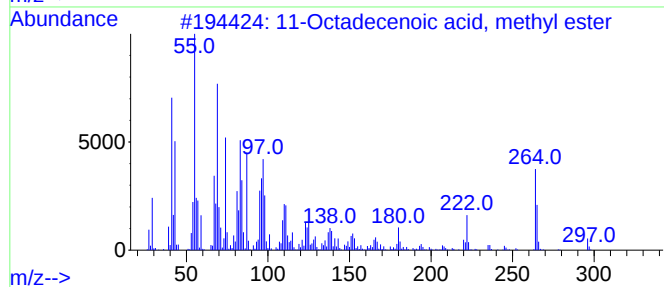
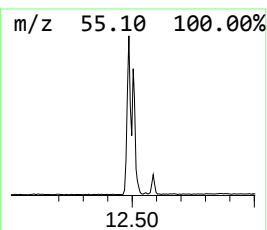
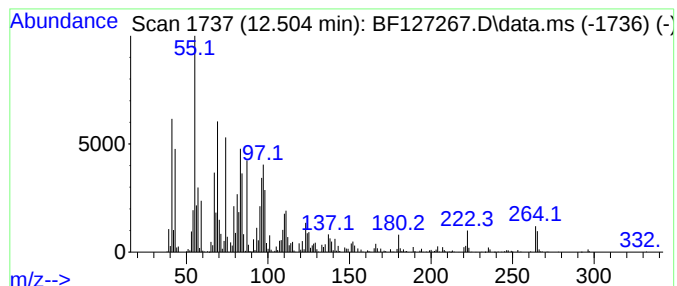
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 11-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.504	58.74 ng	1830230	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
2			6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
3			8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99
4			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
5			9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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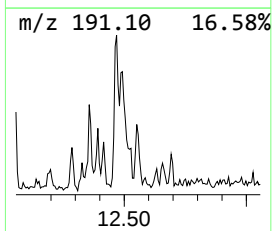
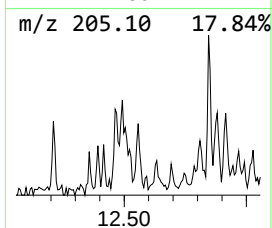
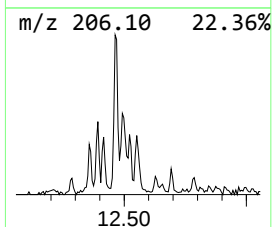
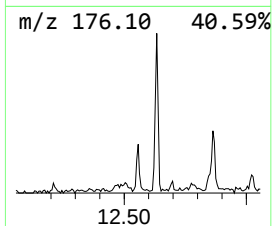
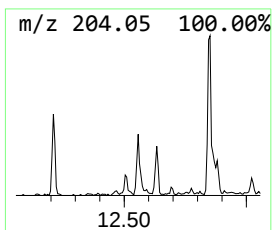
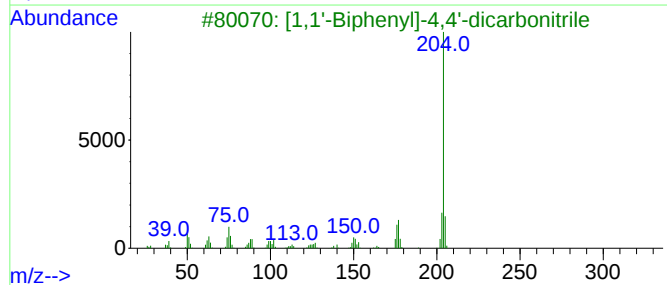
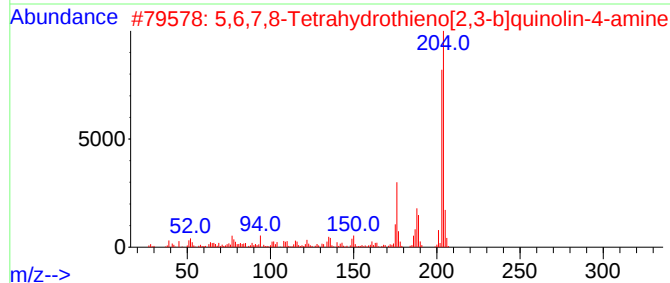
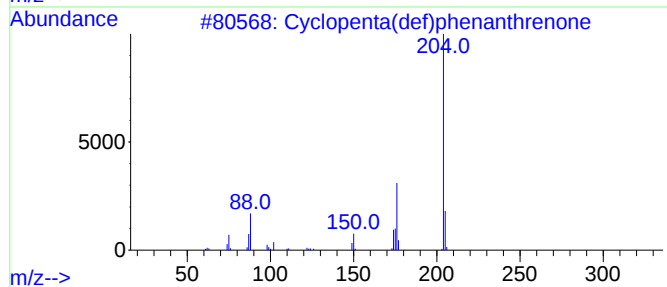
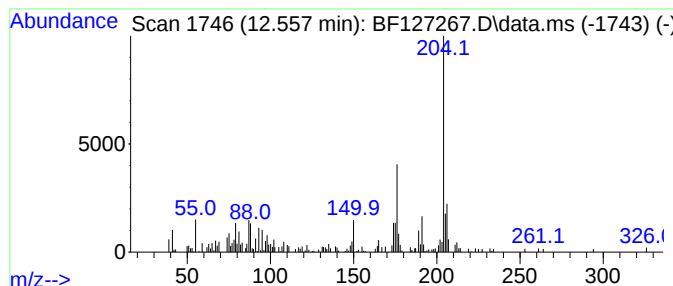
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.557	2.33 ng	72663	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrene	204	C15H8O	005737-13-3	89
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	53
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49
4			2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	43
5			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
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Operator : CG\JU
Sample : N1829-01 5X
Misc :
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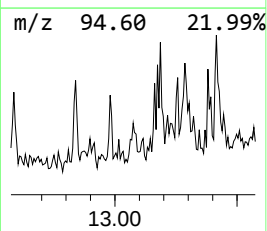
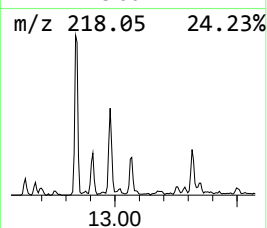
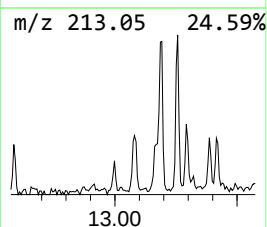
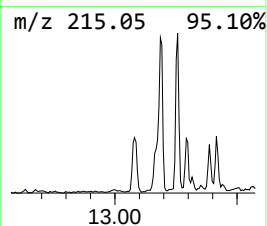
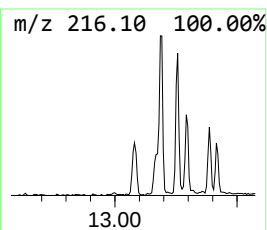
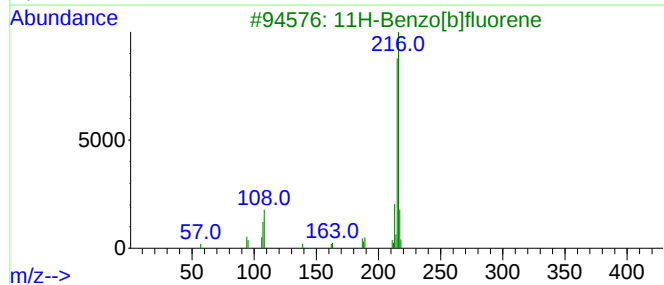
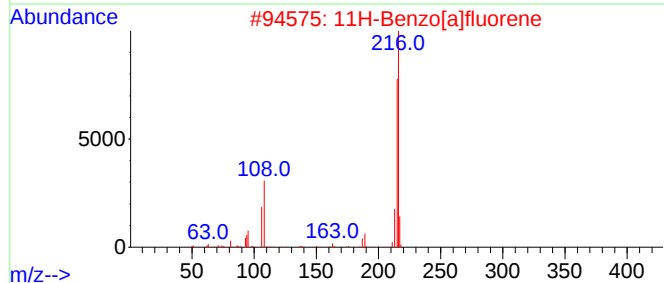
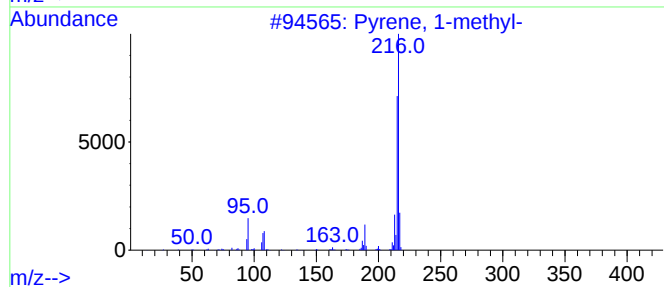
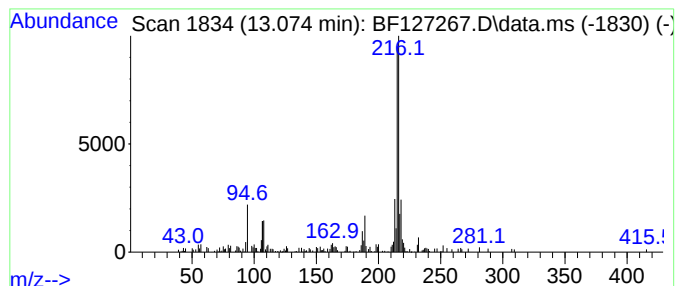
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.074	2.35 ng	108130	Chrysene-d12	14.063	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	89
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	70
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	60
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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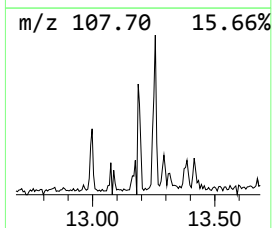
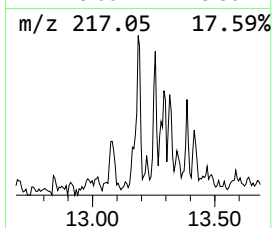
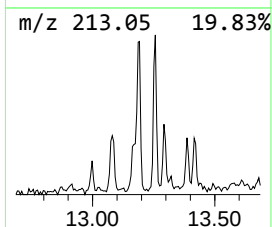
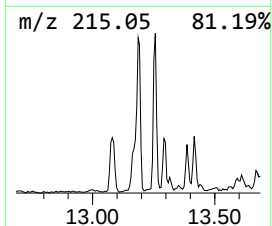
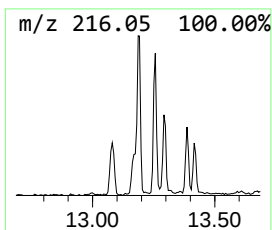
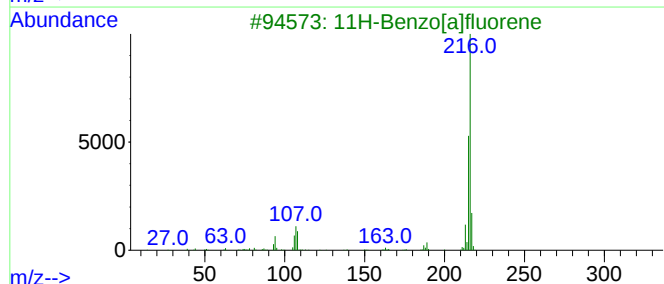
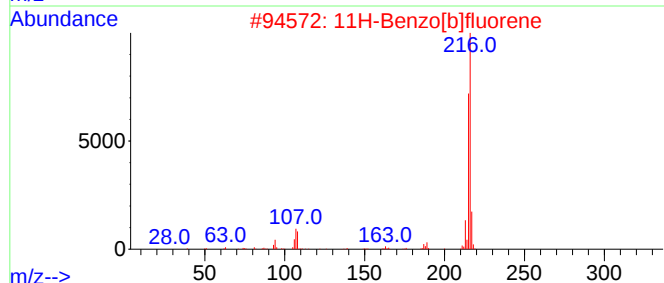
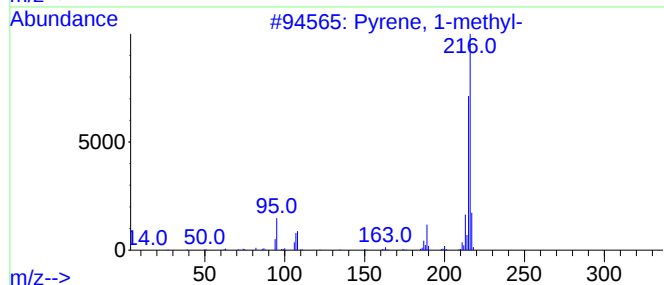
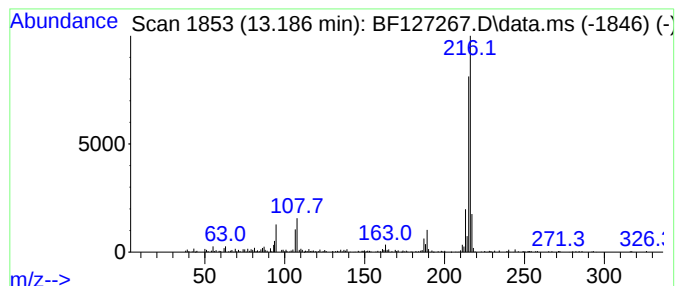
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.186	5.03 ng	231685	Chrysene-d12	14.063

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-030722

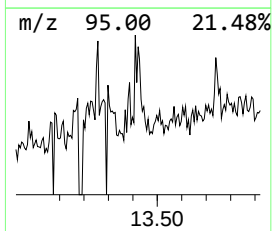
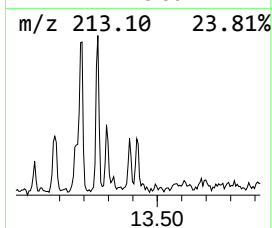
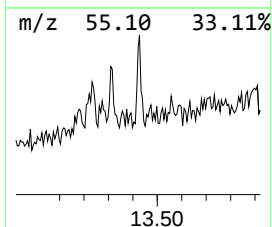
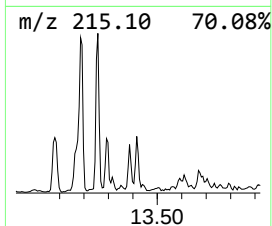
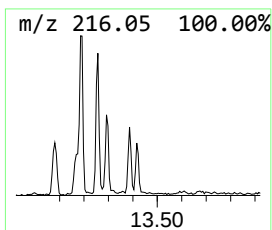
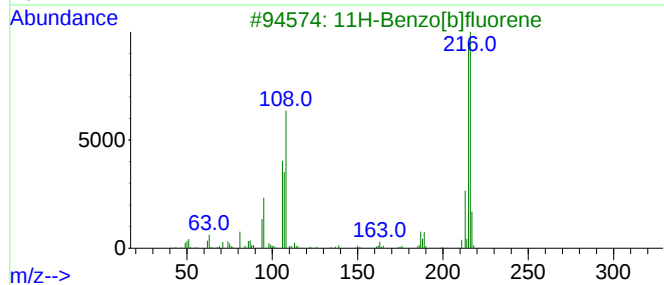
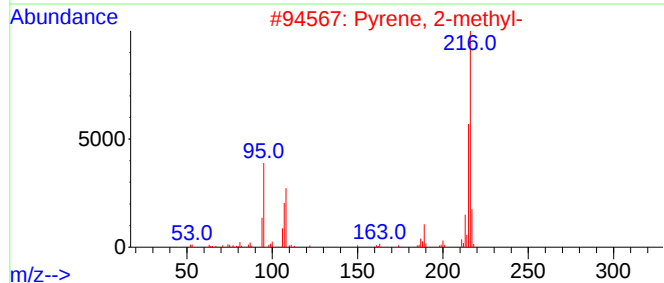
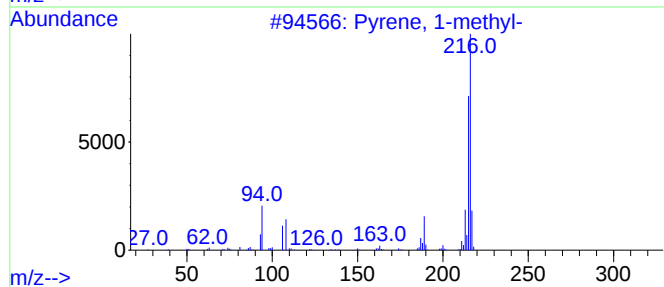
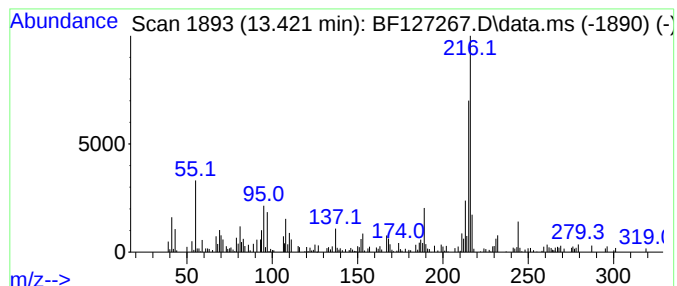
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pyrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.422	2.06 ng	95157	Chrysene-d12	14.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	81
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	70
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
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Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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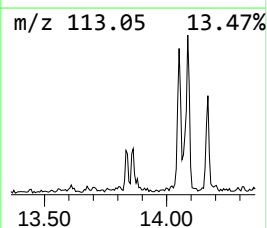
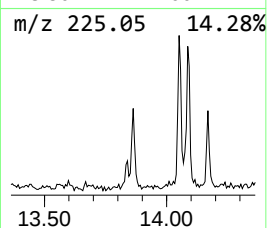
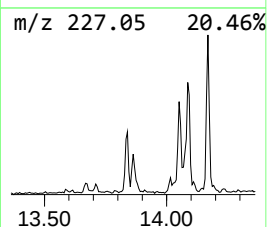
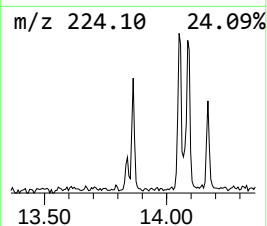
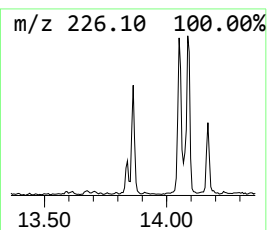
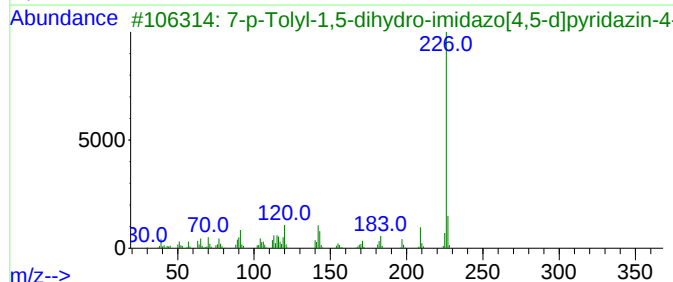
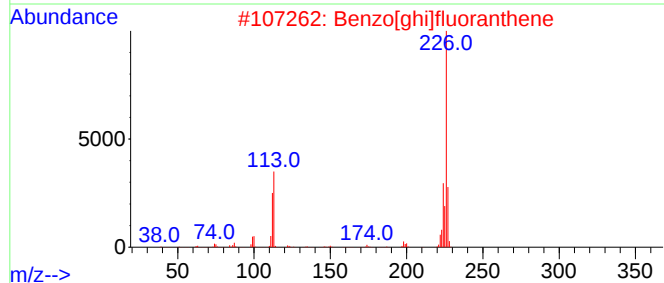
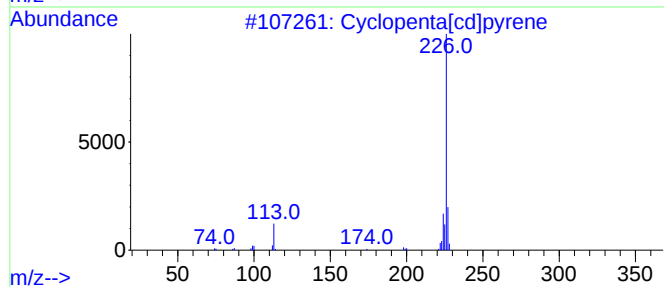
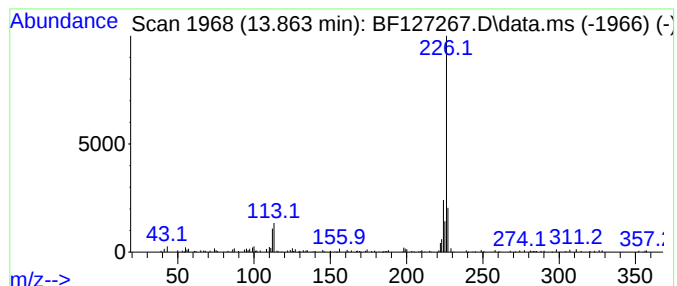
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cyclopenta[cd]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.863	2.02 ng	93328	Chrysene-d12	14.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	93
2			Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	90
3			7-p-Tolyl-1,5-dihydro-imidazo[4,...	226	C12H10N4O	1000317-75-5	53
4			1,3-Diamino-5,6-dihydro-7-methyl...	226	C13H14N4	037436-37-6	53
5			9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	53



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Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-030722

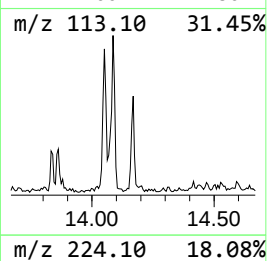
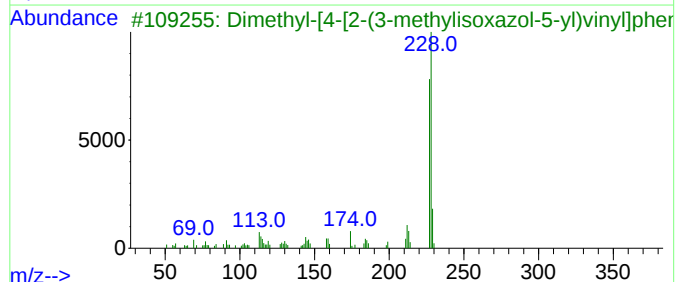
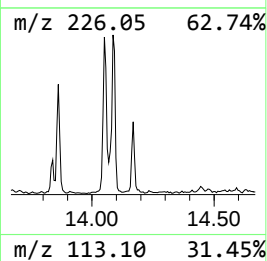
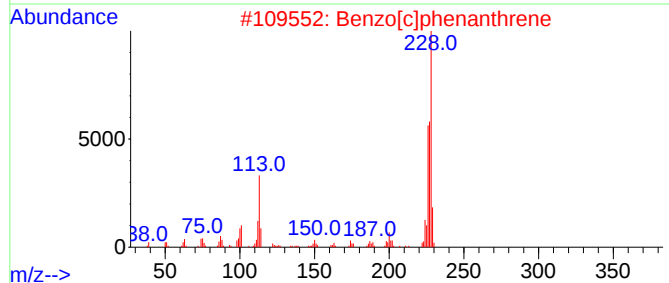
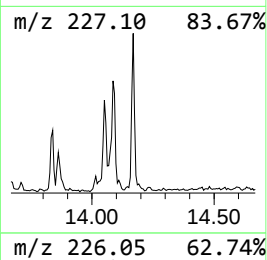
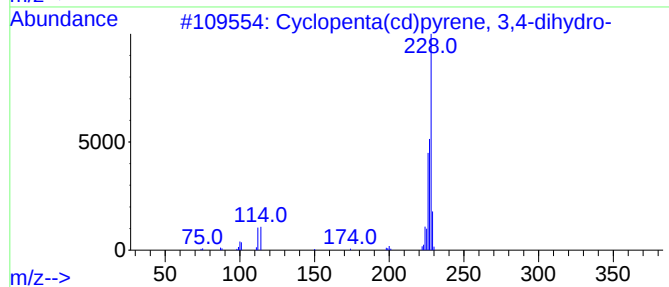
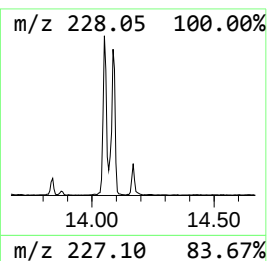
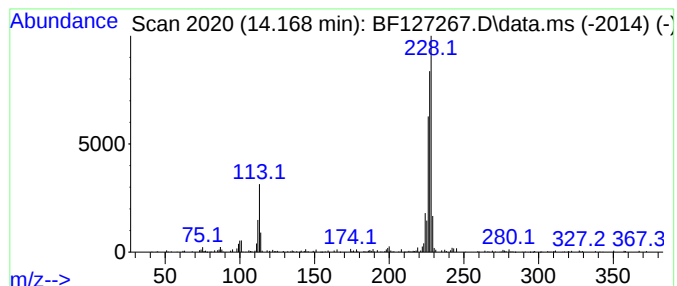
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.169	4.28 ng	197425	Chrysene-d12	14.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	87
2			Benzo[c]phenanthrene	228	C18H12	000195-19-7	58
3			Dimethyl-[4-[2-(3-methylisoxazol...	228	C14H16N2O	1000306-39-6	50
4			5-Phenylthieno[2,3-d]pyrimidin-4-ol	228	C12H8N2OS	035978-39-3	50
5			Diphenylvinylphosphine oxide	228	C14H13OP	002096-78-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-030722

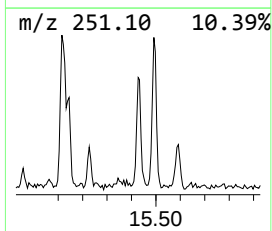
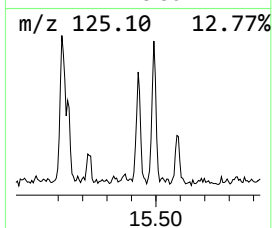
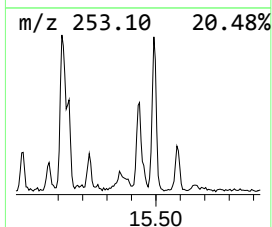
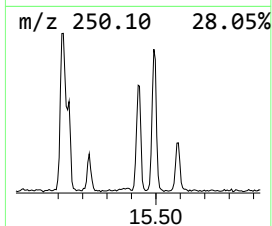
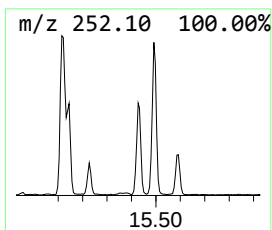
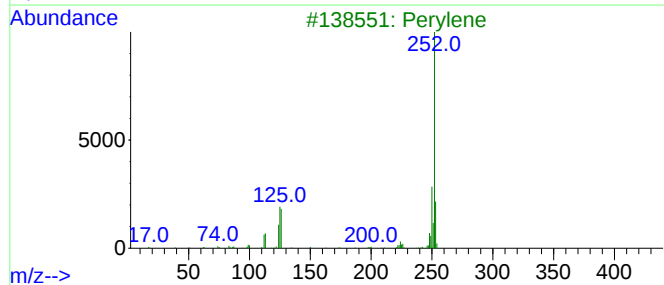
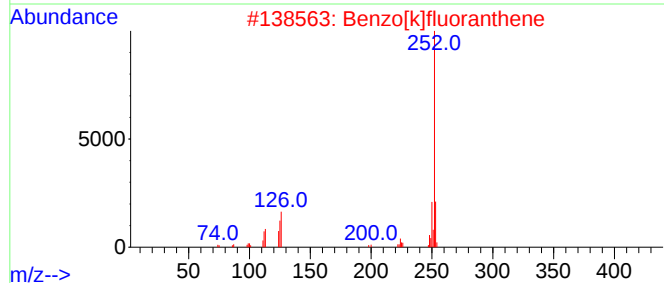
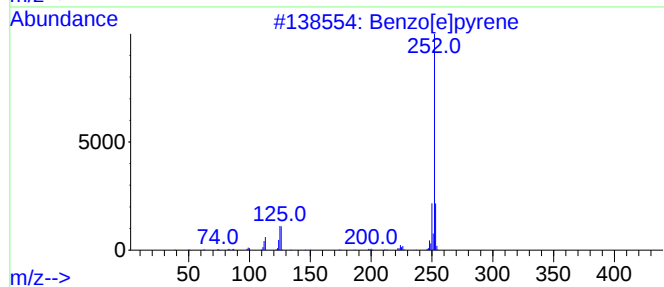
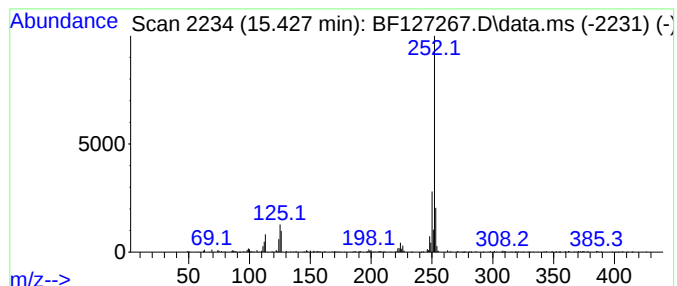
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.427	13.97 ng	231240	Perylene-d12	15.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Perylene	252	C20H12	000198-55-0	94
4			Benzo[a]pyrene	252	C20H12	000050-32-8	94
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030822\
Data File : BF127267.D
Acq On : 08 Mar 2022 19:28
Operator : CG\JU
Sample : N1829-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-030722

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.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
Hexadecanoic ac...	11.792	28.9	ng	900597	4	11.416	623211	20.0
Anthracene, 2-m...	11.945	2.4	ng	74453	4	11.416	623211	20.0
4H-Cyclopenta[d...	12.027	5.9	ng	182910	4	11.416	623211	20.0
Tricyclo[8.2.2...	12.210	2.0	ng	63429	4	11.416	623211	20.0
9,12-Octadecadi...	12.486	116.4	ng	3628260	4	11.416	623211	20.0
11-Octadecenoic...	12.504	58.7	ng	1830230	4	11.416	623211	20.0
Cyclopenta(def)...	12.557	2.3	ng	72663	4	11.416	623211	20.0
Pyrene, 1-methyl-	13.074	2.4	ng	108130	5	14.063	921763	20.0
11H-Benzo[b]flu...	13.186	5.0	ng	231685	5	14.063	921763	20.0
Pyrene, 2-methyl-	13.422	2.1	ng	95157	5	14.063	921763	20.0
Cyclopenta[cd]p...	13.863	2.0	ng	93328	5	14.063	921763	20.0
Cyclopenta(cd)p...	14.169	4.3	ng	197425	5	14.063	921763	20.0
Benzo[e]pyrene	15.427	14.0	ng	231240	6	15.557	331010	20.0