

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
 Data File : BF127109.D
 Acq On : 28 Feb 2022 19:38
 Operator : CG\JU
 Sample : N1681-04 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-1-8FT

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: BF127109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.975	283	287	293	rVB	13292	17242	1.65%	0.108%
2	5.516	545	549	555	rBV	1057636	906412	86.90%	5.658%
3	6.522	716	720	728	rVB	1158349	956175	91.67%	5.968%
4	6.898	781	784	788	rBV	897776	749904	71.89%	4.681%
5	7.463	876	880	883	rBV	780637	641200	61.47%	4.002%
6	8.187	999	1003	1006	rBV	1198441	958416	91.88%	5.982%
7	8.898	1121	1124	1127	rBV	24938	19497	1.87%	0.122%
8	8.998	1138	1141	1143	rVB	28814	22852	2.19%	0.143%
9	9.257	1181	1185	1188	rBV	1395292	1043111	100.00%	6.511%
10	9.357	1200	1202	1206	rBV3	18209	17094	1.64%	0.107%
11	9.528	1227	1231	1234	rBV4	11544	13913	1.33%	0.087%
12	9.598	1239	1243	1245	rBV	28723	30923	2.96%	0.193%
13	9.716	1260	1263	1265	rBV	16842	14424	1.38%	0.090%
14	9.804	1275	1278	1280	rBV	21674	19884	1.91%	0.124%
15	9.857	1284	1287	1291	rVB	14317	15912	1.53%	0.099%
16	9.945	1298	1302	1305	rBV	1002471	901029	86.38%	5.624%
17	9.975	1305	1307	1311	rVB	156774	115720	11.09%	0.722%
18	10.098	1322	1328	1330	rBV4	11421	13694	1.31%	0.085%
19	10.145	1330	1336	1340	rBV	81730	79696	7.64%	0.497%
20	10.339	1365	1369	1370	rBV	16564	18007	1.73%	0.112%
21	10.357	1370	1372	1375	rVB2	19819	21720	2.08%	0.136%
22	10.492	1388	1395	1399	rVB	68586	71405	6.85%	0.446%
23	10.575	1407	1409	1412	rVB3	17233	13831	1.33%	0.086%
24	10.645	1419	1421	1426	rVB	27242	22551	2.16%	0.141%
25	10.704	1428	1431	1432	rBV2	19532	15294	1.47%	0.095%
26	10.728	1432	1435	1439	rVB	575351	505398	48.45%	3.155%
27	10.839	1451	1454	1456	rBV2	21012	23591	2.26%	0.147%
28	11.022	1482	1485	1487	rBV4	16595	21646	2.08%	0.135%
29	11.163	1505	1509	1511	rBV4	13851	16597	1.59%	0.104%
30	11.233	1518	1521	1525	rVB	72827	64854	6.22%	0.405%
31	11.281	1525	1529	1532	rBV3	23176	28679	2.75%	0.179%
32	11.328	1532	1537	1541	rVB	57684	63034	6.04%	0.393%
33	11.375	1541	1545	1548	rBV6	18447	20431	1.96%	0.128%
34	11.433	1551	1555	1557	rBV	895955	741349	71.07%	4.628%
35	11.457	1557	1559	1562	rVV	846904	617731	59.22%	3.856%
36	11.510	1562	1568	1570	rVB	116181	104121	9.98%	0.650%

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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-BF021622.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.592	1579	1582	1583	rBV2	13757	13638	1.31%	0.085%
38	11.663	1590	1594	1599	rBV2	99889	101849	9.76%	0.636%
39	11.728	1603	1605	1608	rVB4	19301	16288	1.56%	0.102%
40	11.763	1608	1611	1615	rBV3	26768	31963	3.06%	0.200%
41	11.810	1617	1619	1623	rVB4	13809	15684	1.50%	0.098%
42	11.886	1629	1632	1635	rBV5	15149	16085	1.54%	0.100%
43	11.933	1637	1640	1643	rVV2	92880	132263	12.68%	0.826%
44	11.963	1643	1645	1648	rVV	125220	106884	10.25%	0.667%
45	12.010	1651	1653	1656	rVB	33537	24872	2.38%	0.155%
46	12.045	1656	1659	1662	rBV2	104437	130685	12.53%	0.816%
47	12.069	1662	1663	1666	rVB	70198	42352	4.06%	0.264%
48	12.116	1669	1671	1673	rBV2	14868	15251	1.46%	0.095%
49	12.228	1687	1690	1692	rBV	57386	49212	4.72%	0.307%
50	12.257	1692	1695	1698	rVB	115228	90446	8.67%	0.565%
51	12.380	1714	1716	1718	rBV2	20189	15955	1.53%	0.100%
52	12.410	1719	1721	1723	rVB	20723	15674	1.50%	0.098%
53	12.486	1730	1734	1737	rBV2	52380	72736	6.97%	0.454%
54	12.569	1745	1748	1753	rBV	54893	61439	5.89%	0.384%
55	12.651	1758	1762	1765	rVV	1077060	868510	83.26%	5.421%
56	12.710	1770	1772	1774	rVV3	18234	15917	1.53%	0.099%
57	12.739	1774	1777	1780	rVB5	22686	26269	2.52%	0.164%
58	12.780	1781	1784	1785	rVV2	16085	17590	1.69%	0.110%
59	12.804	1785	1788	1790	rVB	30271	30670	2.94%	0.191%
60	12.880	1795	1801	1806	rVV	886877	901473	86.42%	5.627%
61	12.927	1806	1809	1812	rVB2	38290	36567	3.51%	0.228%
62	13.016	1821	1824	1833	rVB	841297	785839	75.34%	4.905%
63	13.098	1833	1838	1841	rBV3	56908	101247	9.71%	0.632%
64	13.133	1841	1844	1845	rVV	20447	18400	1.76%	0.115%
65	13.180	1849	1852	1853	rBV2	42706	46357	4.44%	0.289%
66	13.204	1854	1856	1859	rVB2	74406	74046	7.10%	0.462%
67	13.239	1859	1862	1865	rBV5	19121	20979	2.01%	0.131%
68	13.275	1865	1868	1870	rBV	34647	37949	3.64%	0.237%
69	13.310	1870	1874	1877	rVV3	65427	97211	9.32%	0.607%
70	13.339	1877	1879	1881	rVB3	28711	18125	1.74%	0.113%
71	13.404	1888	1890	1893	rVB	35202	25471	2.44%	0.159%
72	13.433	1893	1895	1898	rBV	41357	40101	3.84%	0.250%
73	13.592	1920	1922	1924	rBV3	25610	26223	2.51%	0.164%
74	13.716	1940	1943	1946	rVB	110058	98516	9.44%	0.615%
75	13.827	1958	1962	1965	rBV2	91170	120651	11.57%	0.753%
76	13.857	1965	1967	1969	rVV	72899	58550	5.61%	0.365%

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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-BF021622.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	13.880	1969	1971	1975	rVV2	60976	88404	8.48%	0.552%
78	13.922	1975	1978	1986	rVB2	130808	185883	17.82%	1.160%
79	14.080	2000	2005	2008	rBV2	753140	958846	91.92%	5.985%
80	14.104	2008	2009	2017	rVB	337553	286922	27.51%	1.791%
81	14.186	2021	2023	2025	rBV	49990	38204	3.66%	0.238%
82	15.139	2181	2185	2188	rBV	226221	342042	32.79%	2.135%
83	15.451	2236	2238	2243	rVB	125860	135893	13.03%	0.848%
84	15.516	2245	2249	2253	rBV	164172	210303	20.16%	1.313%
85	15.580	2256	2260	2264	rBV	339732	446704	42.82%	2.788%

Sum of corrected areas: 16020475

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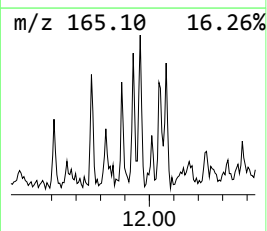
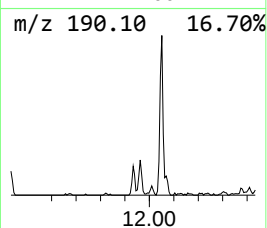
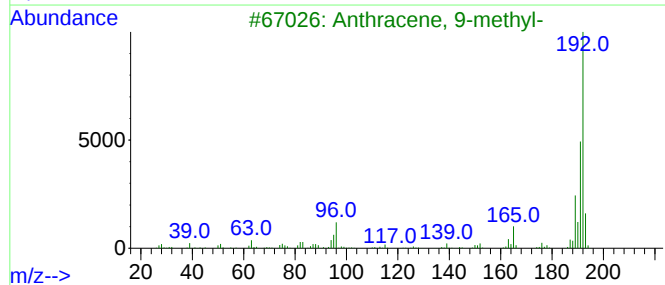
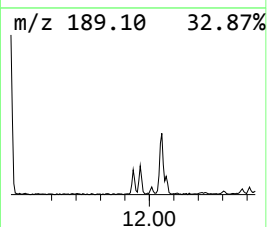
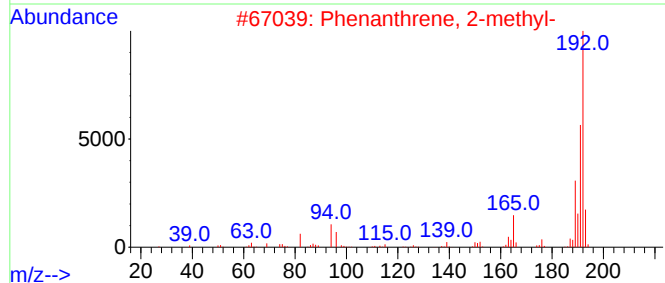
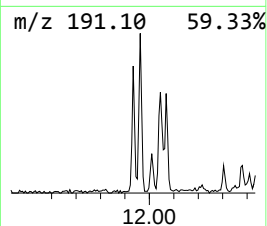
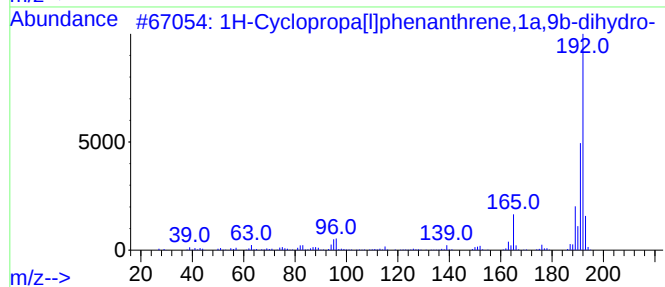
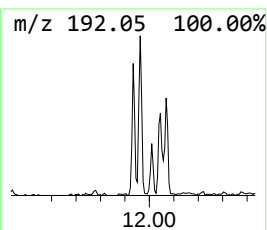
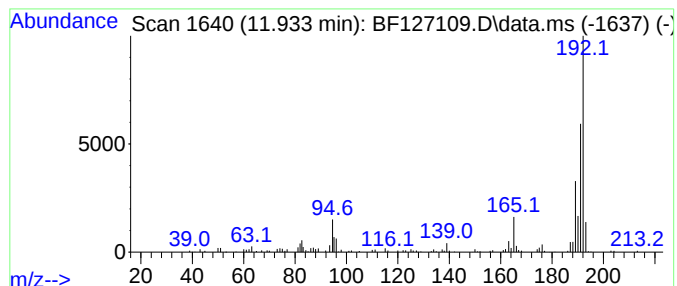
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1H-Cyclopropa[l]phenanthren... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	3.57 ng	132263	Phenanthrene-d10	11.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87



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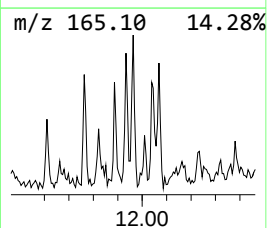
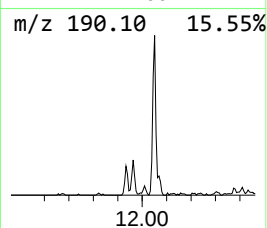
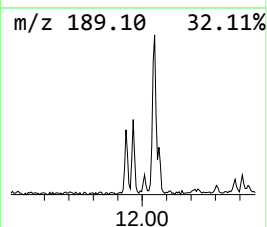
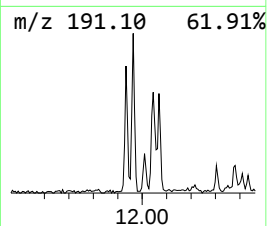
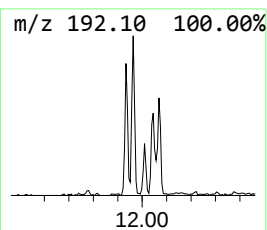
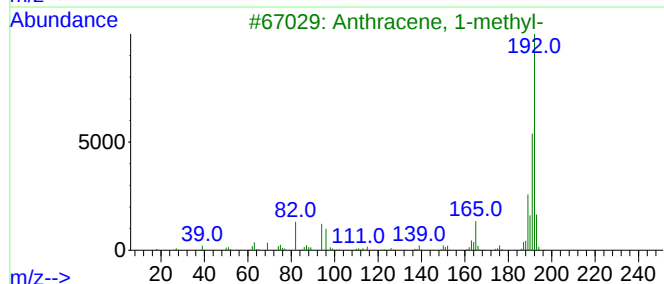
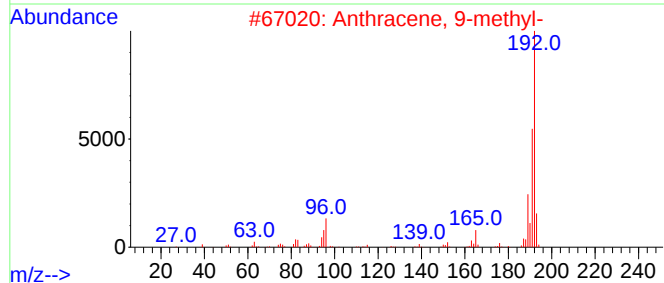
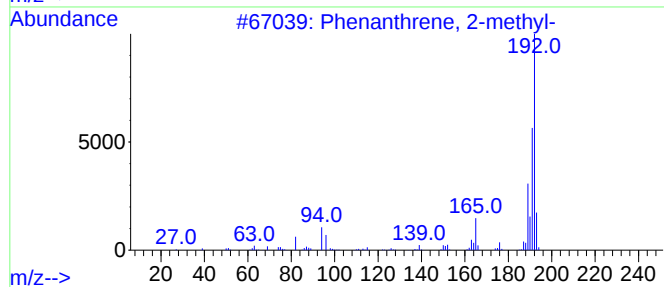
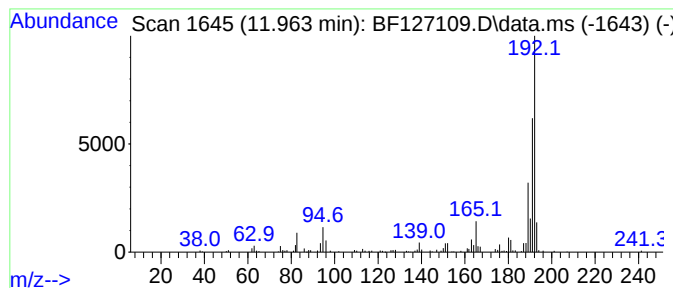
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	2.88 ng	106884	Phenanthrene-d10	11.434

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89



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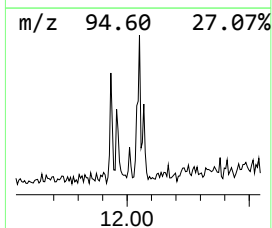
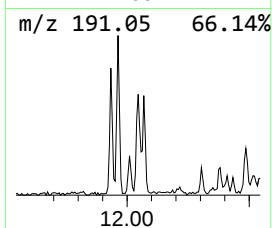
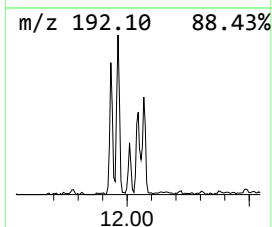
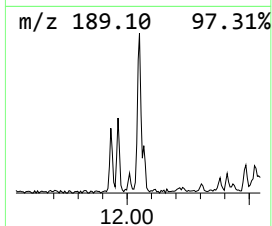
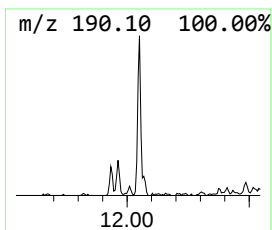
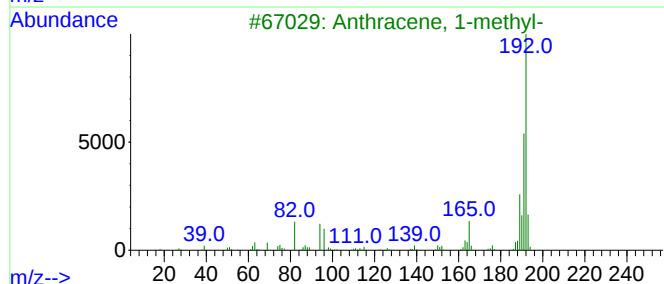
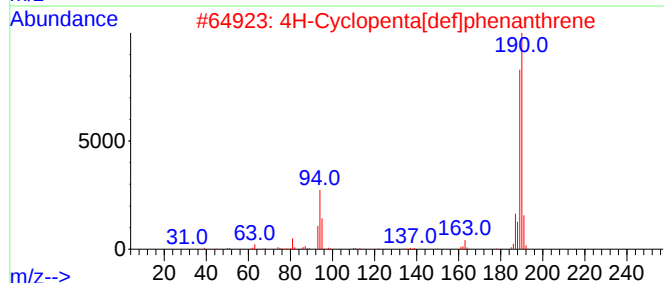
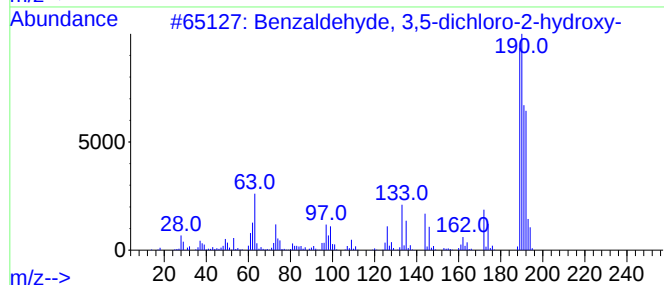
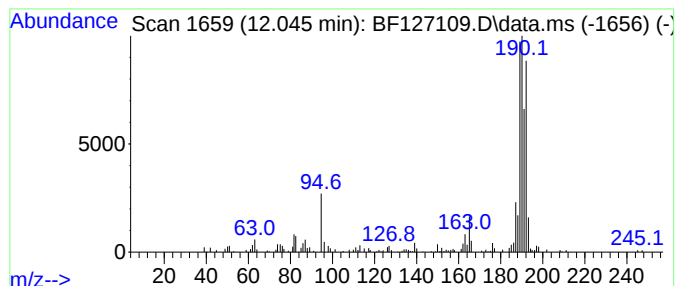
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzaldehyde, 3,5-dichloro-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.045	3.53 ng	130685	Phenanthrene-d10	11.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	41



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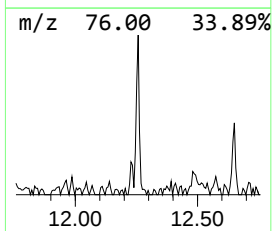
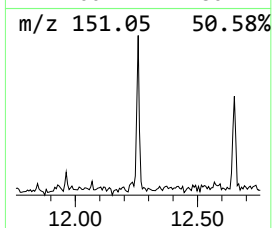
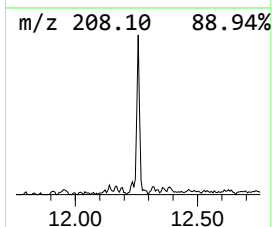
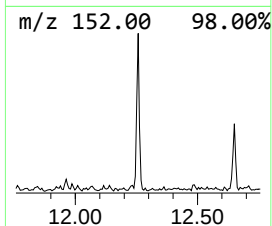
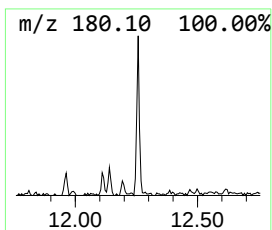
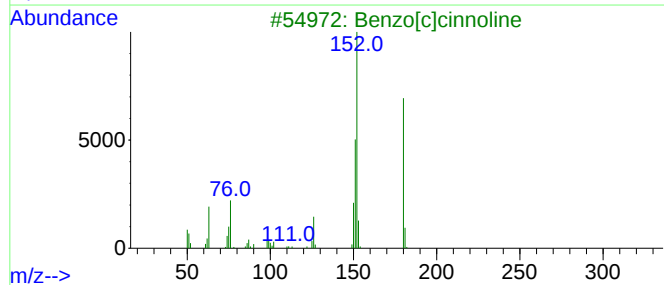
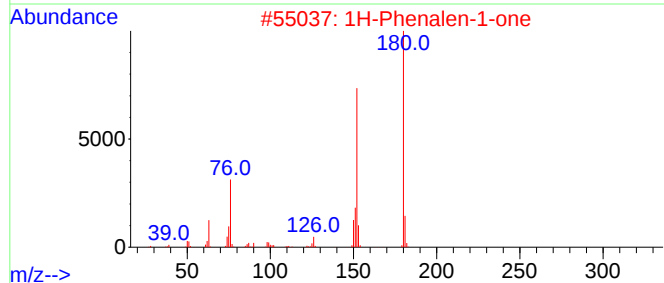
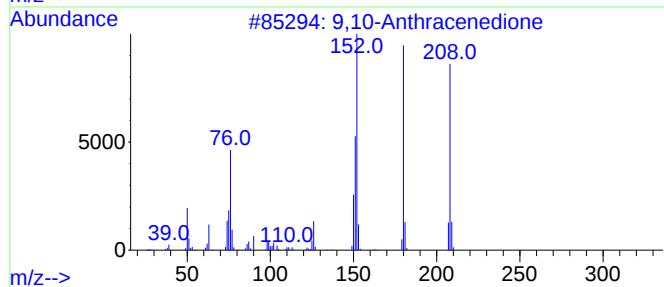
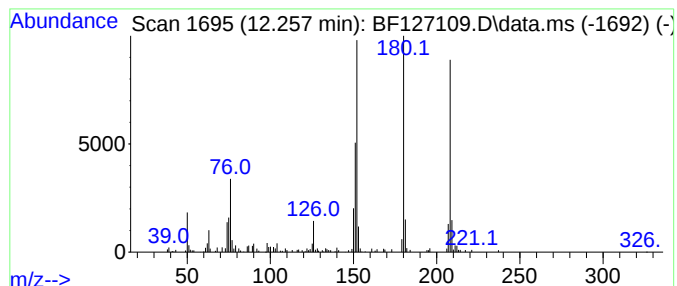
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.257	2.44 ng	90446	Phenanthrene-d10	11.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2			1H-Phenalen-1-one	180	C13H8O	000548-39-0	91
3			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	70
4			9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	60
5			9H-Fluoren-9-one	180	C13H8O	000486-25-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
Sample : N1681-04 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2-1-8FT

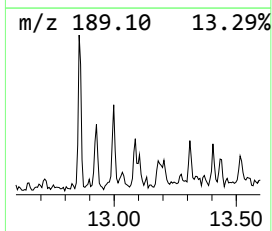
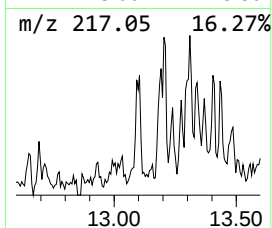
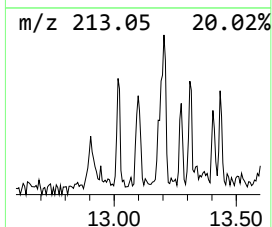
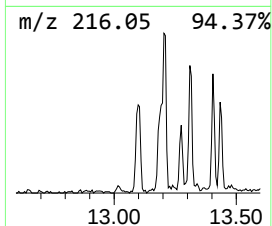
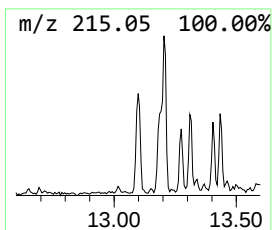
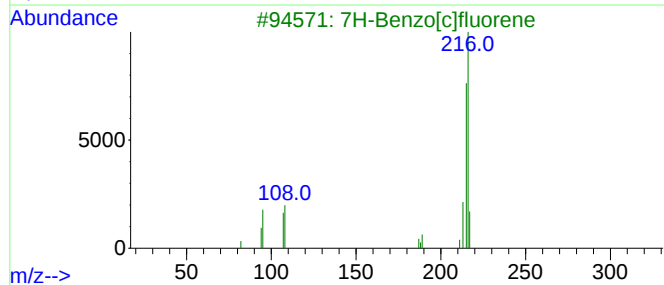
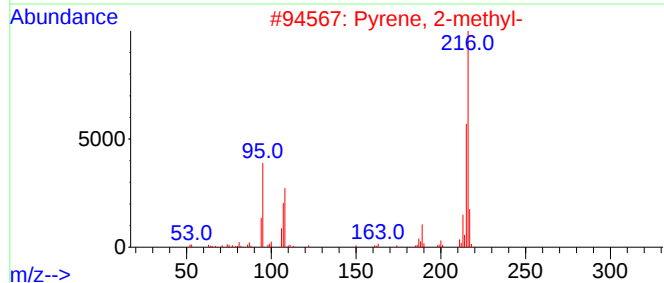
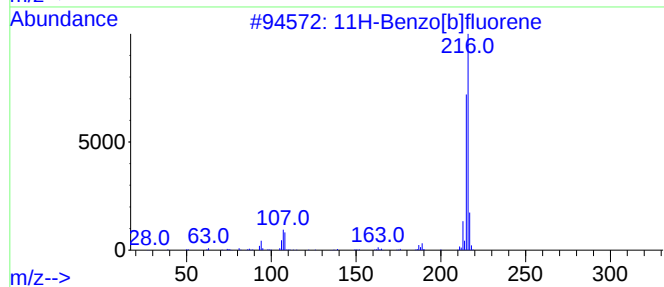
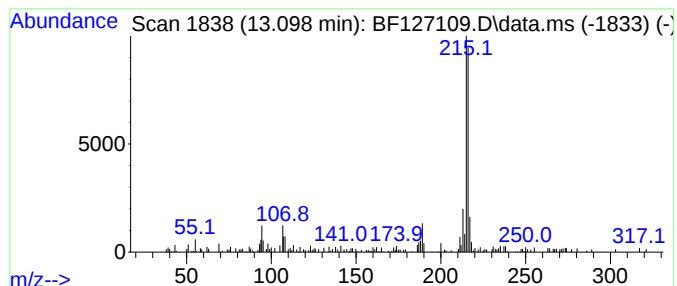
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.098	2.11 ng	101247	Chrysene-d12	14.080

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
3			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	80
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
Sample : N1681-04 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-2-1-8FT

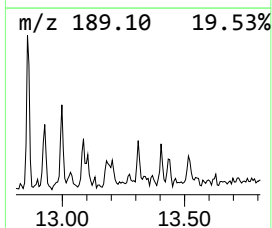
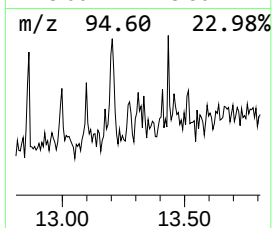
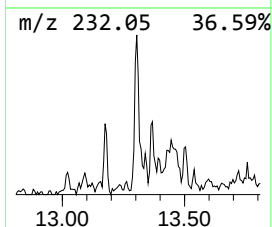
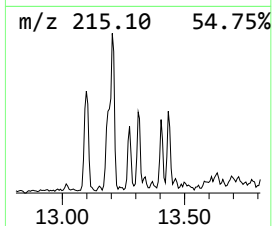
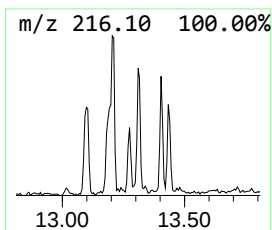
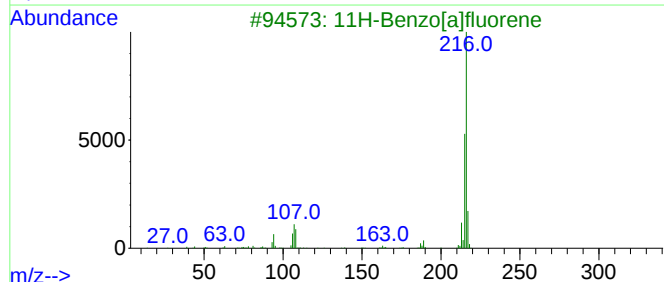
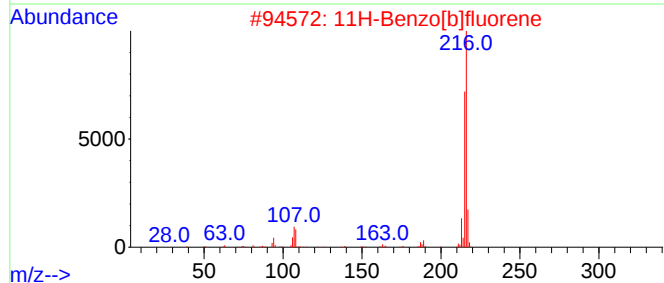
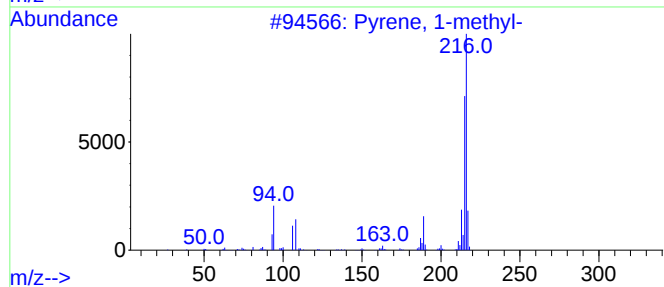
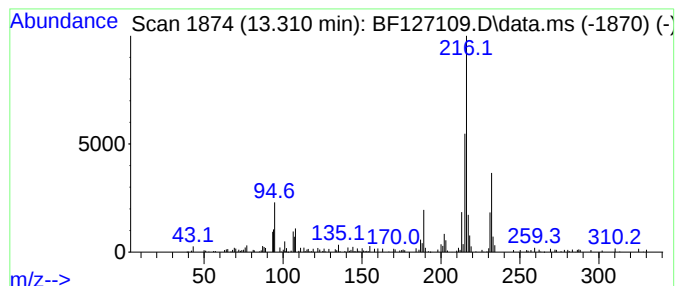
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.310	2.03 ng	97211	Chrysene-d12	14.080

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	64
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	62
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	60
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
Sample : N1681-04 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

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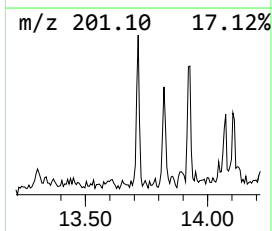
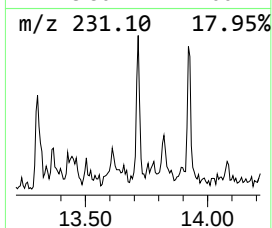
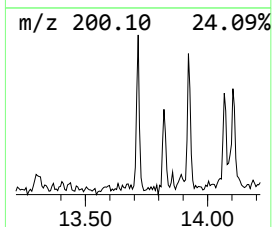
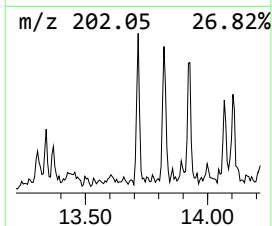
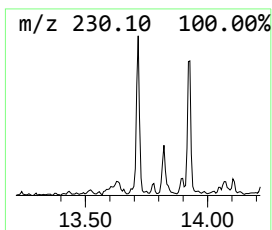
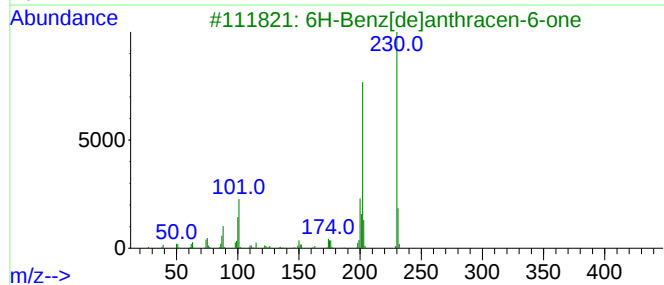
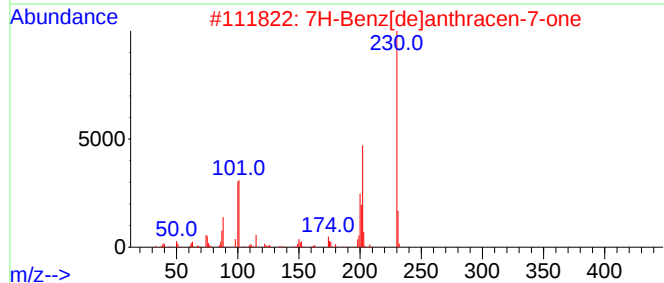
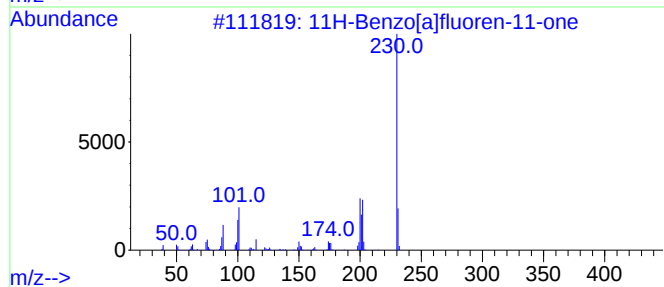
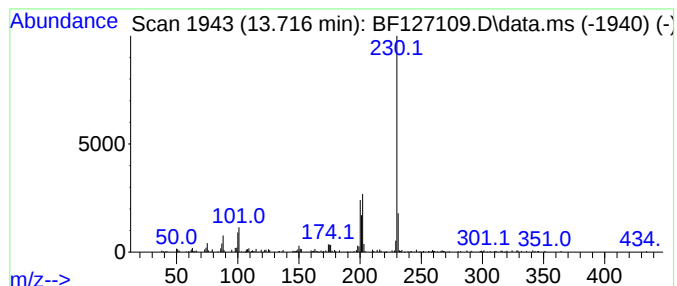
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 11H-Benzo[a]fluoren-11-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	2.05 ng	98516	Chrysene-d12	14.080

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	91
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
3			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	58
4			o-Terphenyl	230	C18H14	000084-15-1	53
5			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
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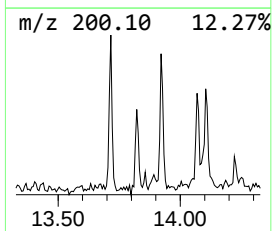
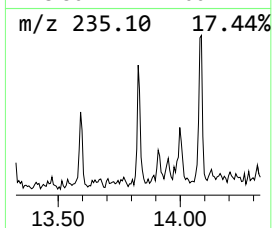
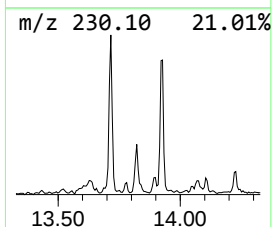
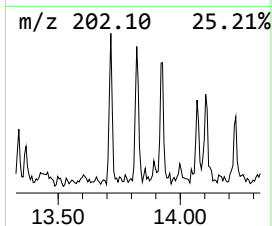
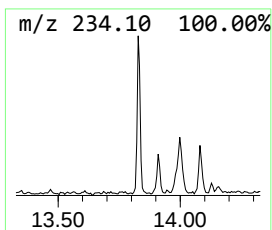
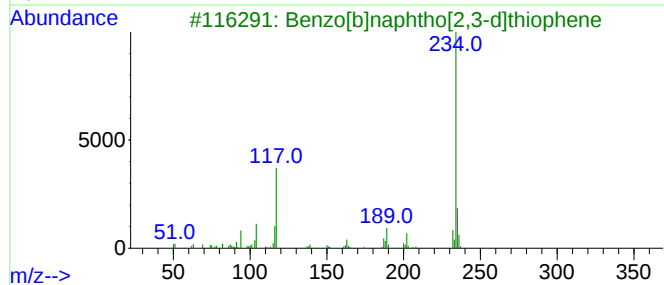
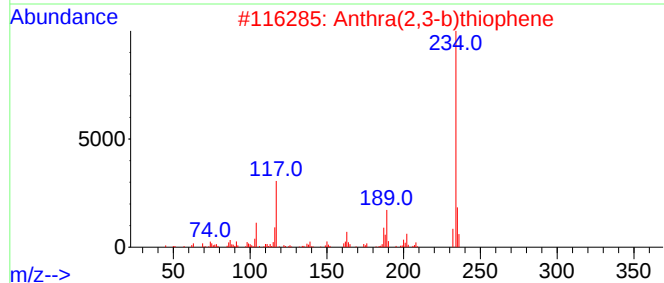
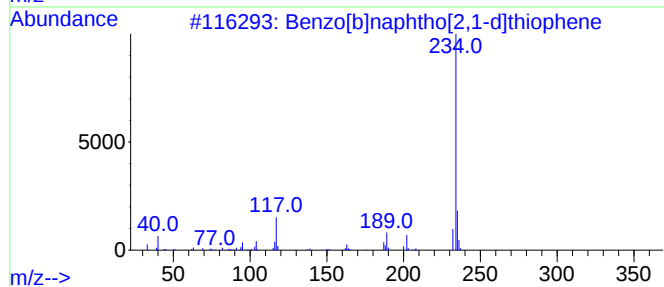
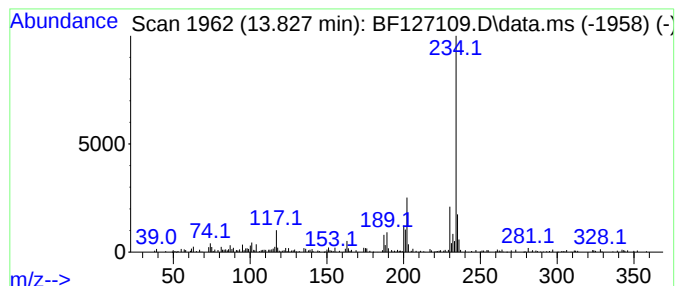
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.827	2.52 ng	120651	Chrysene-d12	14.080

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	83
2			Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	83
3			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	58
4			Phenanthro[4,3-b]thiophene	234	C16H10S	000195-68-6	53
5			Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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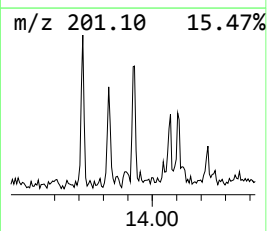
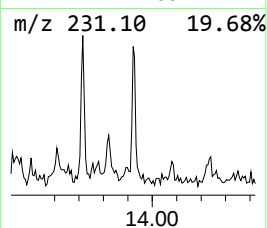
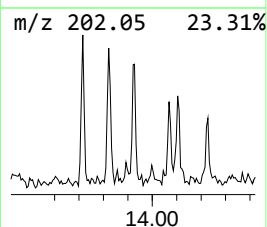
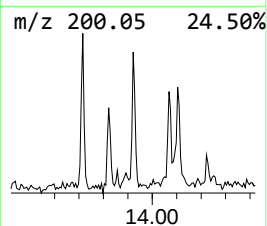
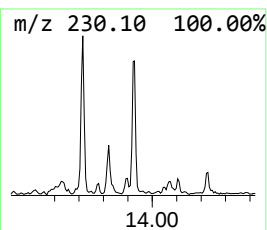
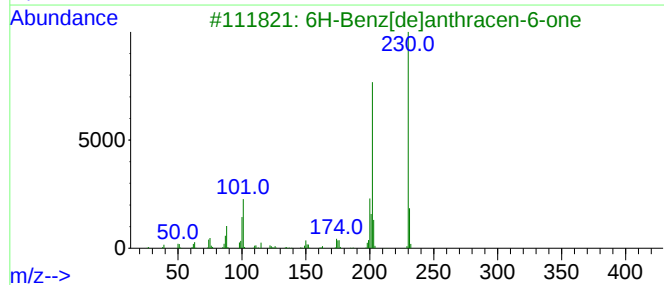
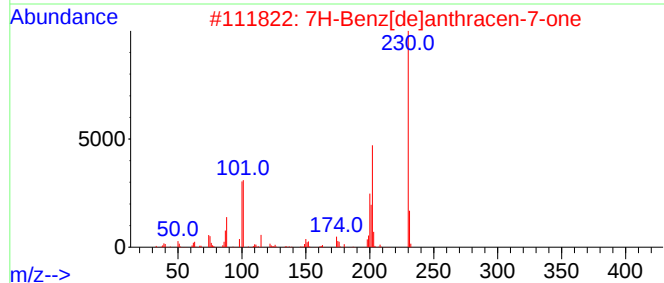
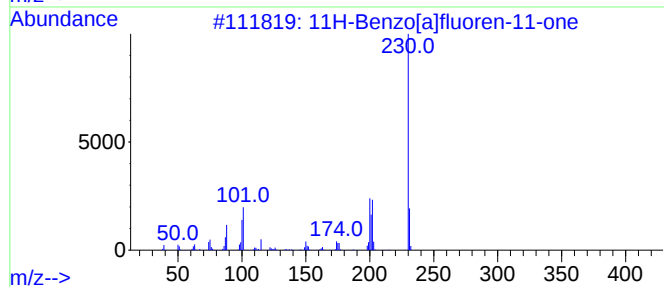
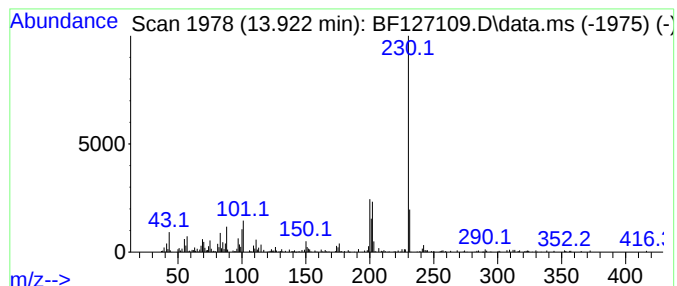
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7H-Benz[de]anthracen-7-one Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.922	3.88 ng	185883	Chrysene-d12	14.080

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
3			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	64
4			o-Terphenyl	230	C18H14	000084-15-1	59
5			.delta.(sup1),.alpha.-Acenaphthe...	230	C15H6N2O	014619-86-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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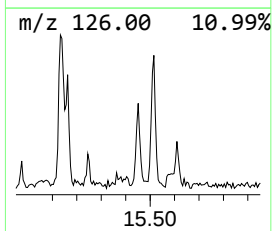
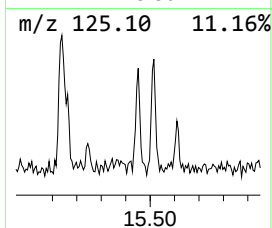
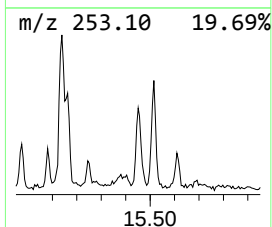
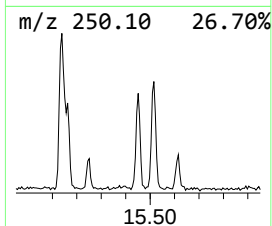
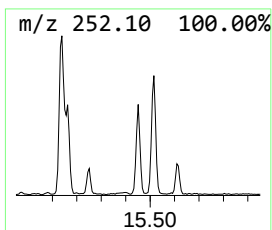
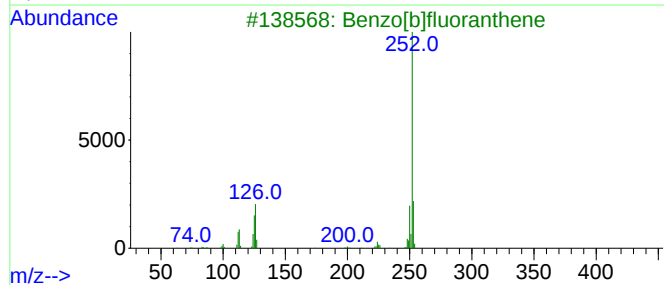
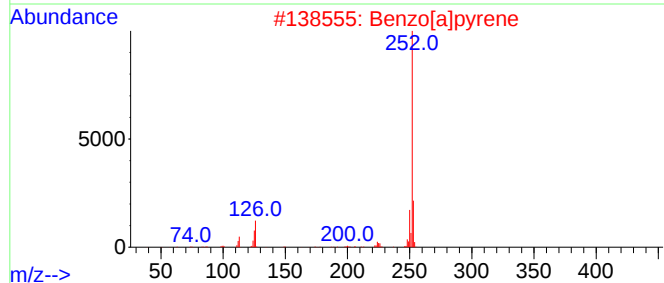
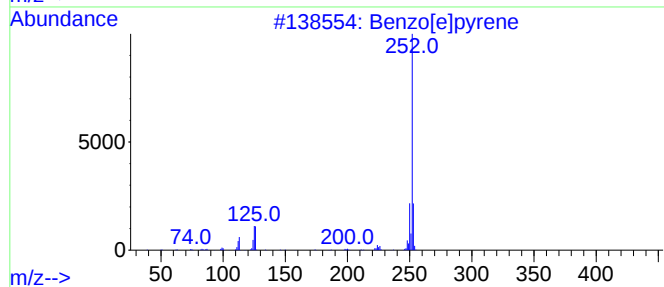
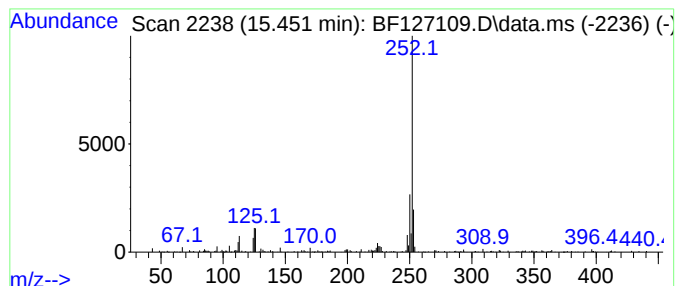
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.451	6.08 ng	135893	Perylene-d12	15.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
5			Perylene	252	C20H12	000198-55-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022822\
Data File : BF127109.D
Acq On : 28 Feb 2022 19:38
Operator : CG\JU
Sample : N1681-04 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2-1-8FT

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
1H-Cyclopropa[1...	11.933	3.6	ng	132263	4	11.434	741349	20.0
Phenanthrene, 2...	11.963	2.9	ng	106884	4	11.434	741349	20.0
Benzaldehyde, 3...	12.045	3.5	ng	130685	4	11.434	741349	20.0
9,10-Anthracene...	12.257	2.4	ng	90446	4	11.434	741349	20.0
11H-Benzo[b]flu...	13.098	2.1	ng	101247	5	14.080	958846	20.0
Pyrene, 1-methyl-	13.310	2.0	ng	97211	5	14.080	958846	20.0
11H-Benzo[a]flu...	13.716	2.0	ng	98516	5	14.080	958846	20.0
Benzo[b]naphtho...	13.827	2.5	ng	120651	5	14.080	958846	20.0
7H-Benz[de]anth...	13.922	3.9	ng	185883	5	14.080	958846	20.0
Benzo[e]pyrene	15.451	6.1	ng	135893	6	15.580	446704	20.0