

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
 Data File : BF141391.D
 Acq On : 04 Feb 2025 01:16
 Operator : RC/JU
 Sample : Q1216-19 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JPP-26.2-012825

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-BF012925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141391.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.987	482	492	500	rBV	24918	38807	3.54%	0.271%
2	5.416	560	565	582	rVB	596204	803416	73.27%	5.605%
3	6.434	733	738	755	rBV	601397	837558	76.38%	5.843%
4	6.792	794	799	805	rBV	413178	511095	46.61%	3.566%
5	7.351	885	894	899	rBV	423867	543627	49.58%	3.793%
6	7.616	933	939	947	rVB2	15789	24742	2.26%	0.173%
7	8.075	1008	1017	1025	rBV	511408	693208	63.22%	4.836%
8	8.369	1060	1067	1073	rBV6	6612	15727	1.43%	0.110%
9	8.498	1085	1089	1095	rBV	12196	16639	1.52%	0.116%
10	8.669	1113	1118	1122	rBV	32889	39531	3.61%	0.276%
11	8.786	1135	1138	1143	rVB	13668	16826	1.53%	0.117%
12	8.886	1151	1155	1160	rVB	13252	17884	1.63%	0.125%
13	9.151	1195	1200	1205	rBV	728557	898936	81.98%	6.272%
14	9.233	1209	1214	1222	rVB2	31000	45734	4.17%	0.319%
15	9.416	1240	1245	1250	rBV	8173	12253	1.12%	0.085%
16	9.486	1252	1257	1260	rBV	16786	24041	2.19%	0.168%
17	9.610	1274	1278	1283	rVB	7910	12280	1.12%	0.086%
18	9.686	1287	1291	1296	rBV	10936	14419	1.31%	0.101%
19	9.763	1296	1304	1308	rBV	9993	15178	1.38%	0.106%
20	9.828	1308	1315	1319	rBV	497188	637344	58.12%	4.446%
21	9.863	1319	1321	1325	rVB	57374	58198	5.31%	0.406%
22	9.986	1338	1342	1345	rBV2	7688	11270	1.03%	0.079%
23	10.033	1345	1350	1354	rBV	31108	38355	3.50%	0.268%
24	10.075	1354	1357	1364	rVB3	8556	12880	1.17%	0.090%
25	10.145	1365	1369	1372	rBV	7924	11042	1.01%	0.077%
26	10.239	1380	1385	1386	rBV3	9358	13451	1.23%	0.094%
27	10.257	1386	1388	1393	rVB3	10656	13417	1.22%	0.094%
28	10.375	1404	1408	1413	rVB	39558	50986	4.65%	0.356%
29	10.445	1413	1420	1422	rBV2	10047	21170	1.93%	0.148%
30	10.539	1431	1436	1441	rVB2	57717	80111	7.31%	0.559%
31	10.616	1444	1449	1455	rBV	281637	366680	33.44%	2.558%
32	10.739	1466	1470	1479	rVB4	19847	34710	3.17%	0.242%
33	10.822	1479	1484	1487	rBV6	6497	11814	1.08%	0.082%
34	10.922	1495	1501	1505	rBV3	16703	36749	3.35%	0.256%
35	11.057	1519	1524	1525	rBV4	15172	21011	1.92%	0.147%
36	11.122	1531	1535	1539	rVB	21619	26970	2.46%	0.188%

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37	11.180	1539	1545	1548	rBV5	26968	52873	4.82%	0.369%
38	11.210	1548	1550	1559	rVB	29766	39956	3.64%	0.279%
39	11.316	1563	1568	1570	rBV	417179	572937	52.25%	3.997%
40	11.339	1570	1572	1577	rVV	512157	573673	52.32%	4.002%
41	11.392	1577	1581	1585	rVV	114826	153370	13.99%	1.070%
42	11.427	1585	1587	1591	rVB2	12213	13653	1.25%	0.095%
43	11.551	1603	1608	1615	rBV2	27628	53237	4.85%	0.371%
44	11.616	1615	1619	1622	rVV	27150	35859	3.27%	0.250%
45	11.651	1622	1625	1630	rVB3	12603	16886	1.54%	0.118%
46	11.816	1649	1653	1655	rVV	63767	81977	7.48%	0.572%
47	11.845	1655	1658	1663	rVV	125041	171738	15.66%	1.198%
48	11.892	1663	1666	1669	rVV	39642	55342	5.05%	0.386%
49	11.927	1669	1672	1681	rVB2	103691	203367	18.55%	1.419%
50	12.022	1682	1688	1691	rBV5	15935	31001	2.83%	0.216%
51	12.116	1699	1704	1707	rBV	49282	78035	7.12%	0.544%
52	12.263	1724	1729	1733	rBV	53201	82065	7.48%	0.573%
53	12.369	1743	1747	1750	rBV	38665	53758	4.90%	0.375%
54	12.410	1750	1754	1758	rVV2	30182	53903	4.92%	0.376%
55	12.451	1758	1761	1766	rVB	41566	57243	5.22%	0.399%
56	12.527	1769	1774	1779	rBV	742595	968552	88.33%	6.757%
57	12.574	1779	1782	1788	rVB3	44485	68846	6.28%	0.480%
58	12.757	1806	1813	1819	rBV	729958	1096559	100.00%	7.650%
59	12.904	1830	1838	1844	rBV	530044	764798	69.75%	5.336%
60	12.974	1845	1850	1854	rBV3	56613	108152	9.86%	0.755%
61	13.039	1858	1861	1863	rBV	43069	54122	4.94%	0.378%
62	13.186	1883	1886	1899	rVB3	100819	216803	19.77%	1.513%
63	13.280	1900	1902	1905	rBV	27343	26989	2.46%	0.188%
64	13.951	2011	2016	2020	rBV2	590801	1059448	96.62%	7.391%
65	14.986	2188	2192	2201	rVB	253184	529683	48.30%	3.695%
66	15.280	2239	2242	2248	rVB	148165	236429	21.56%	1.649%
67	15.345	2249	2253	2258	rVB	191075	271873	24.79%	1.897%
68	15.404	2259	2263	2267	rBV	203903	338983	30.91%	2.365%
69	16.839	2503	2507	2517	rVB2	81833	183479	16.73%	1.280%

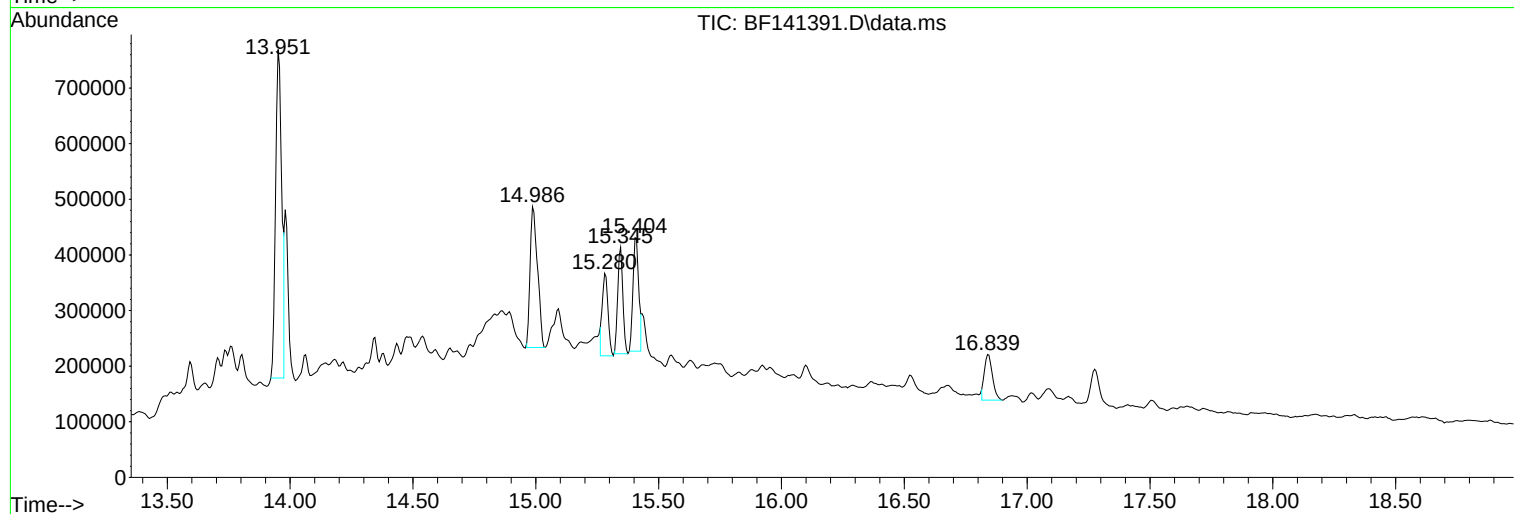
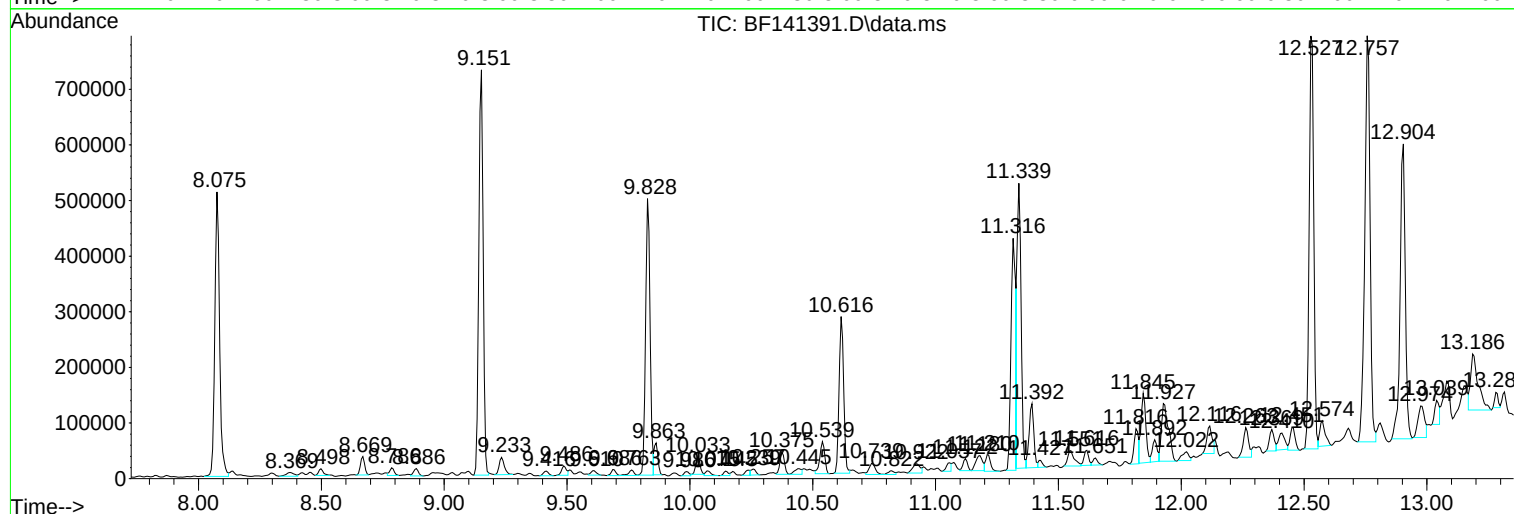
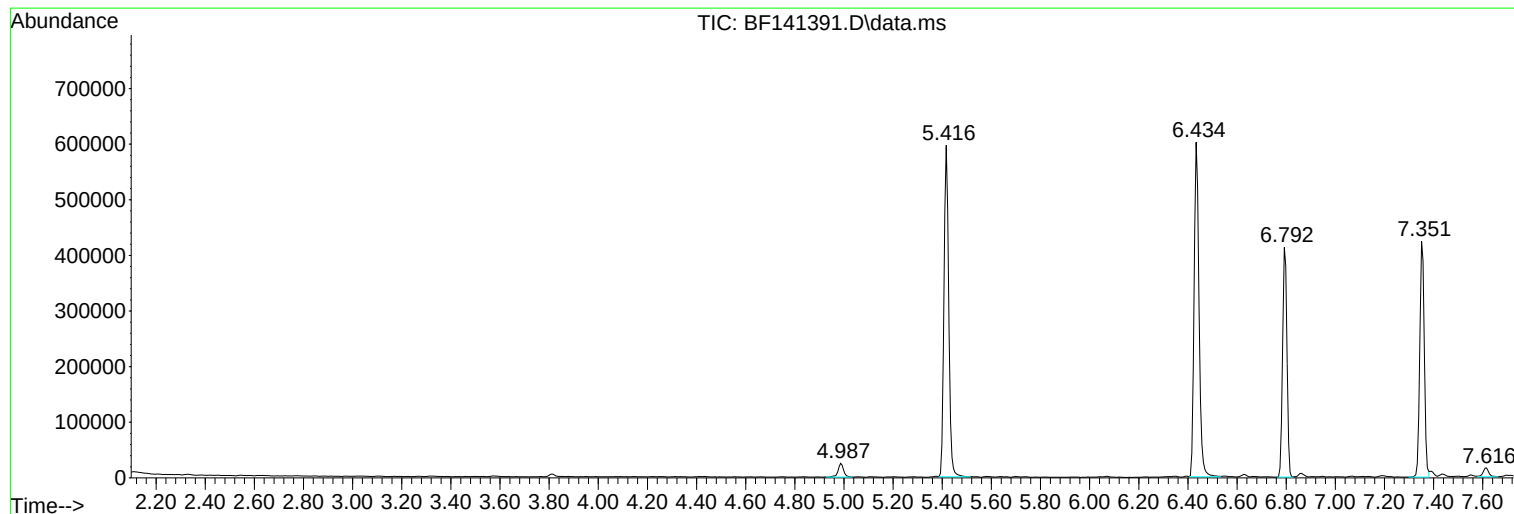
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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JPP-26.2-012825

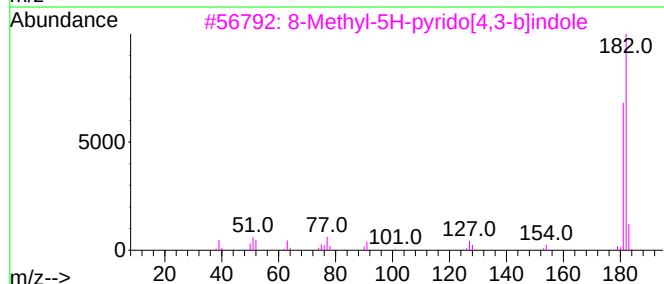
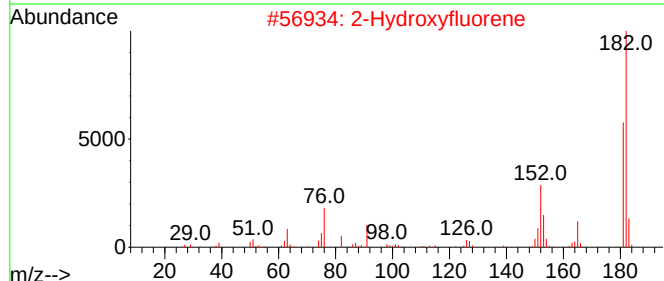
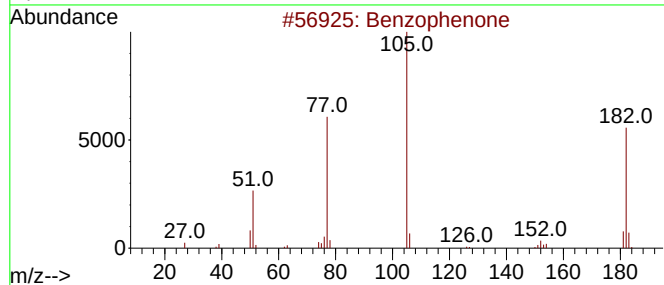
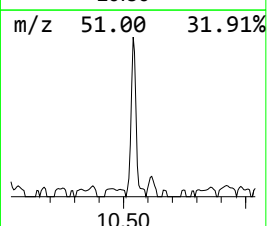
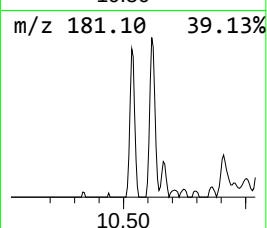
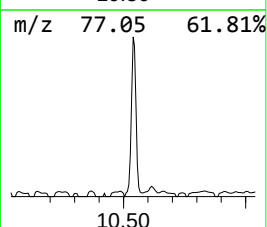
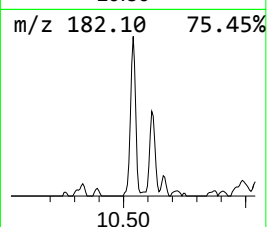
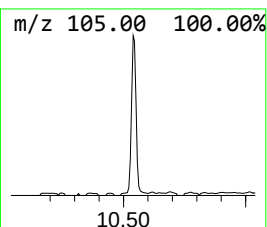
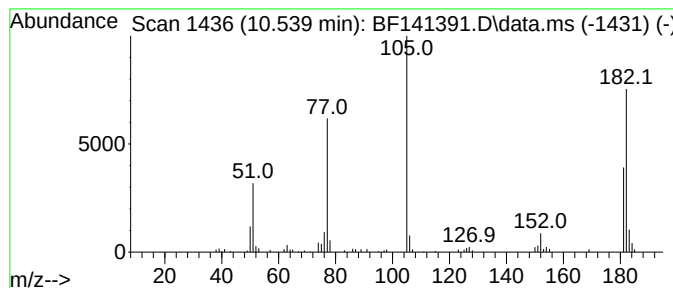
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.539	2.51 ng	80111	Acenaphthene-d10	9.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	87
2			2-Hydroxyfluorene	182	C13H10O	002443-58-5	49
3			8-Methyl-5H-pyrido[4,3-b]indole	182	C12H10N2	088894-13-7	49
4			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	30
5			Benzhydrazide, N2-(2-methoxy-5-n...	299	C15H13N3O4	349572-84-5	27



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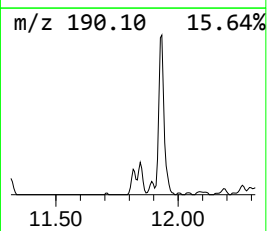
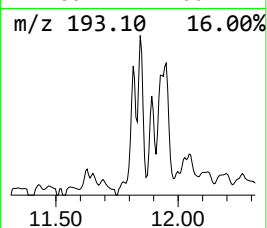
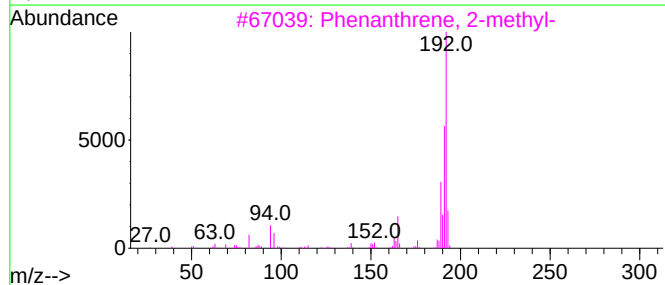
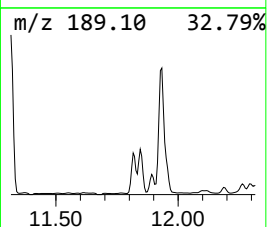
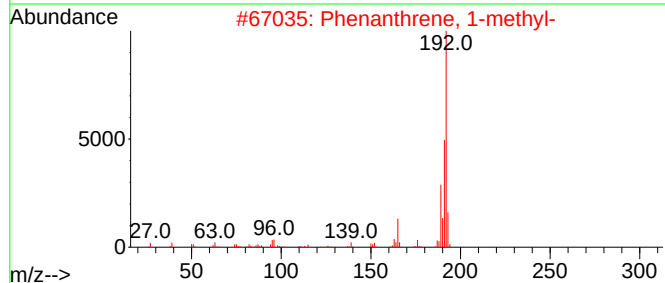
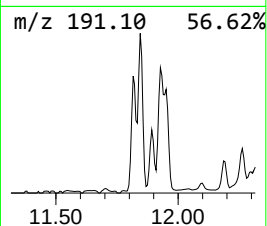
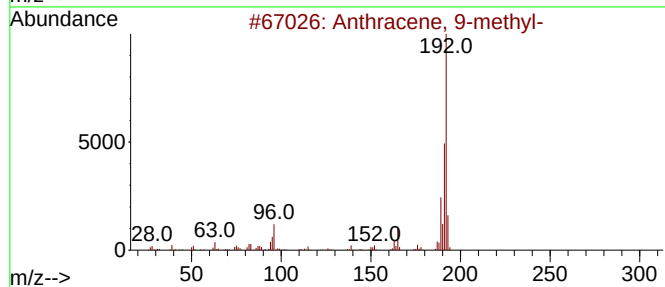
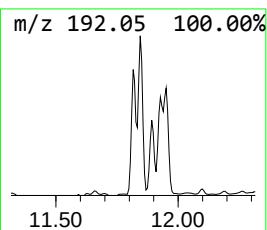
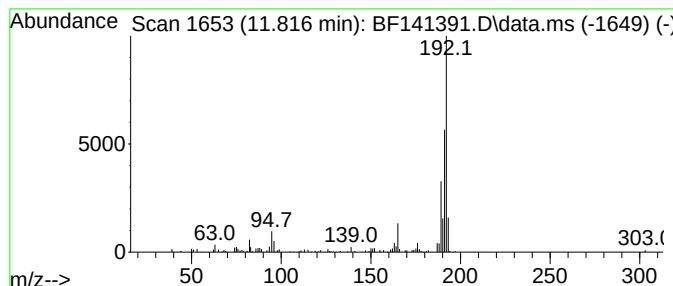
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Anthracene, 9-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.816	2.86 ng	81977	Phenanthrene-d10	11.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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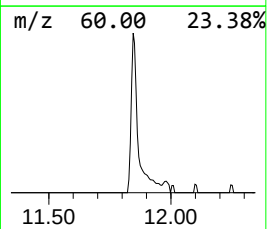
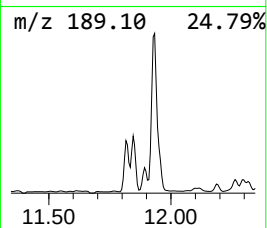
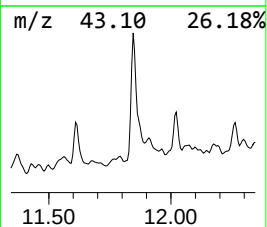
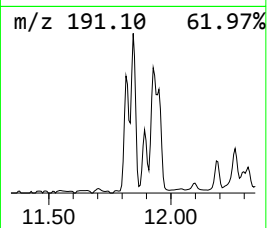
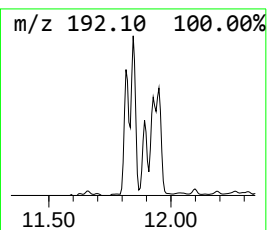
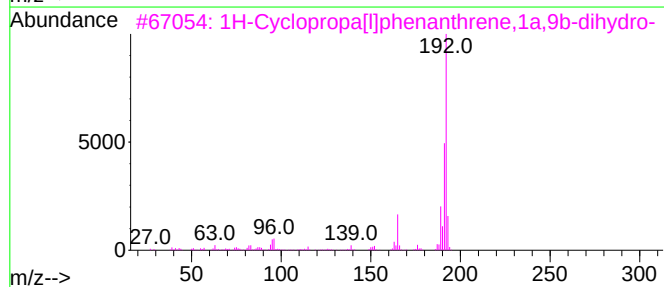
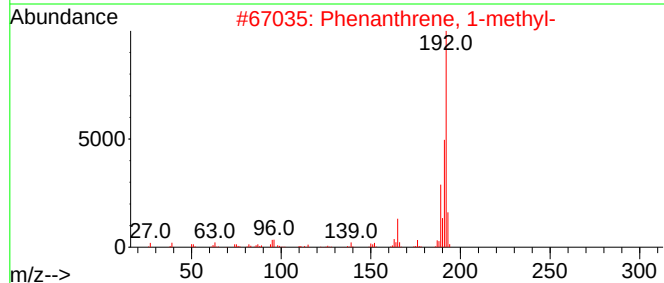
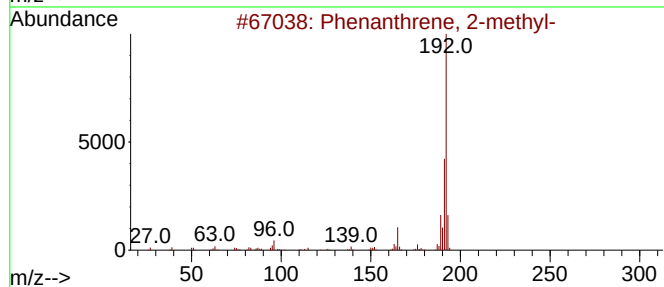
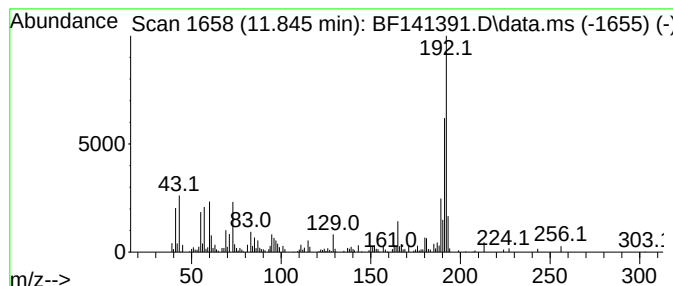
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Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.845	6.00 ng	171738	Phenanthrene-d10	11.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93



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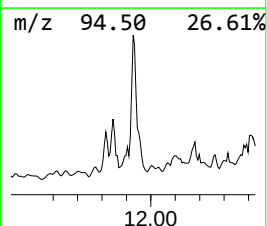
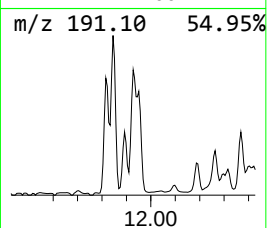
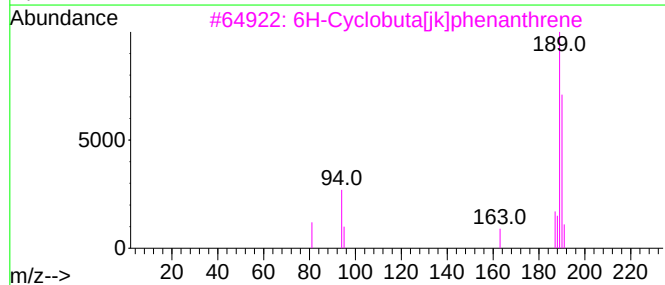
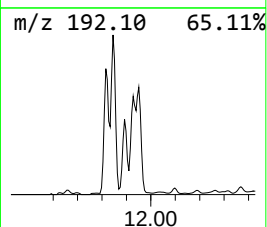
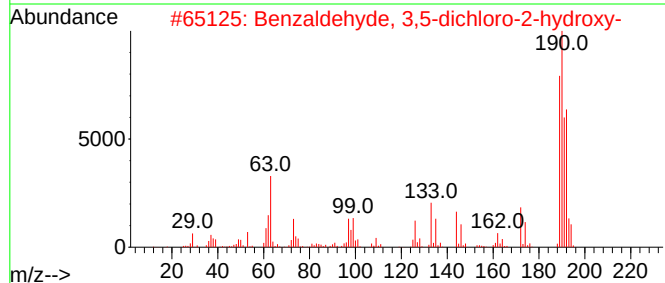
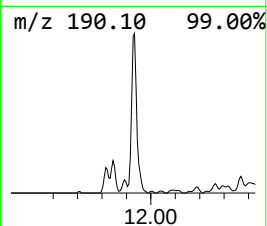
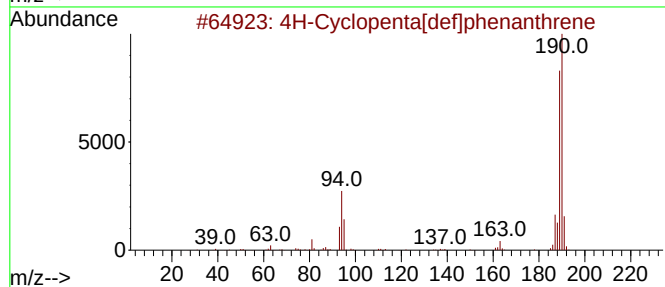
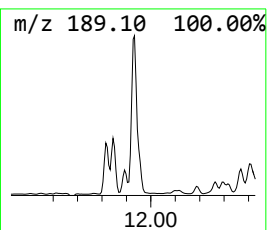
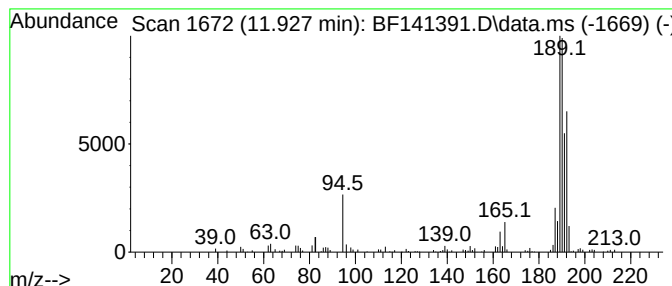
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.927	7.10 ng	203367	Phenanthrene-d10	11.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	25
5			Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JPP-26.2-012825

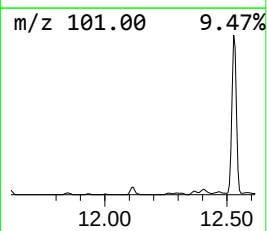
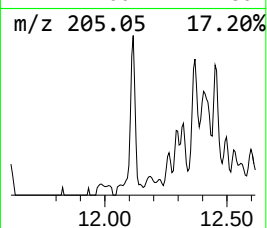
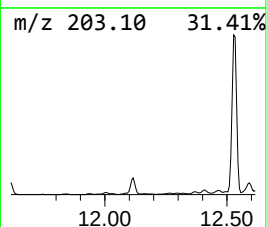
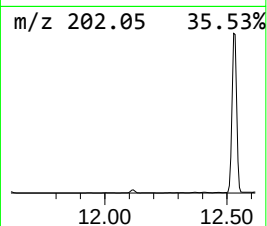
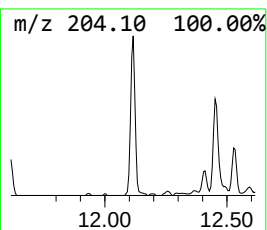
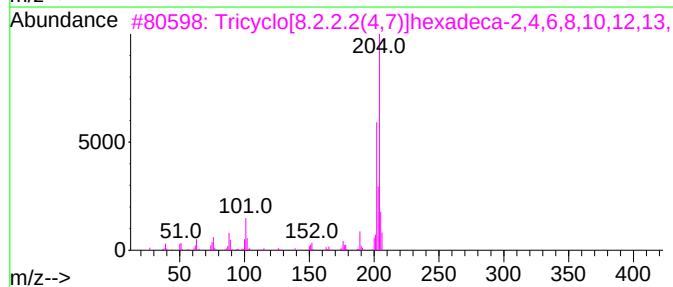
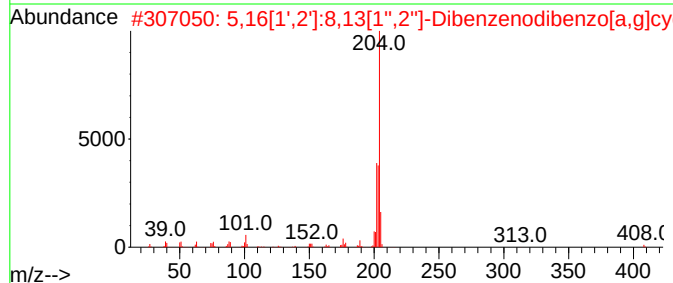
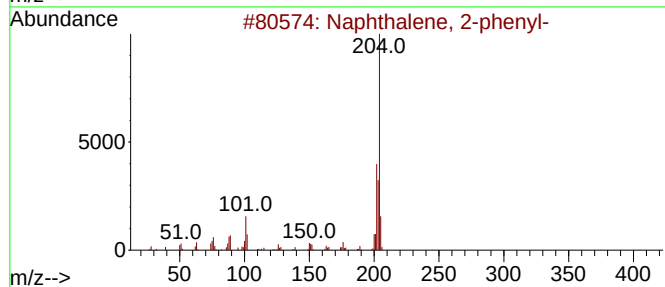
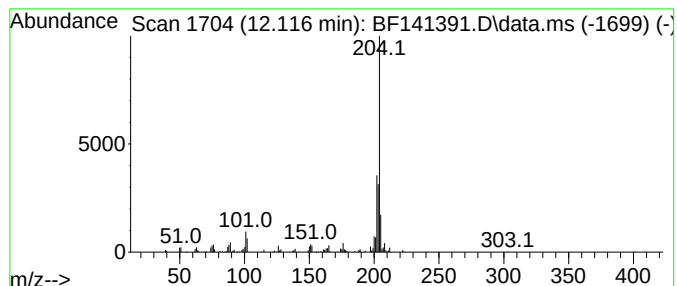
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.116	2.72 ng	78035	Phenanthrene-d10	11.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	95
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	74
4			9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	62
5			9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
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Instrument :
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ClientSampleId :
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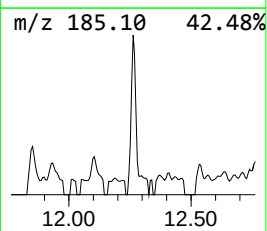
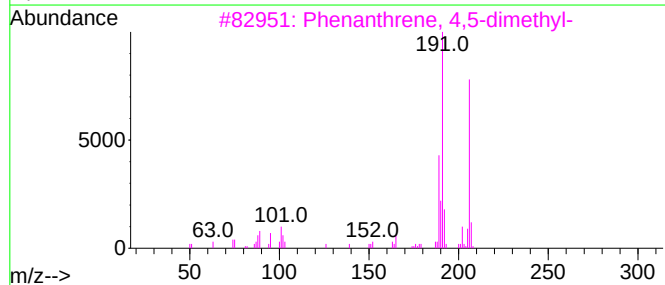
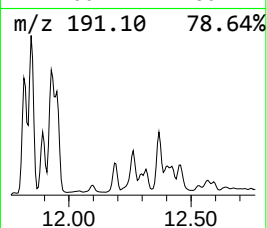
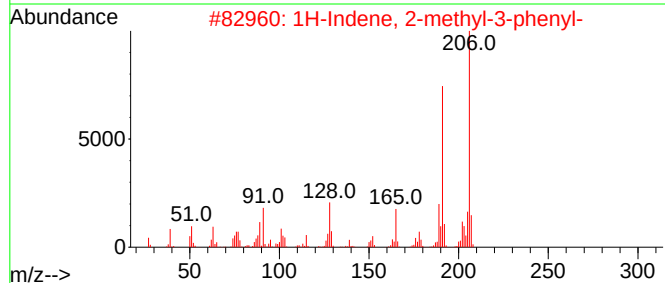
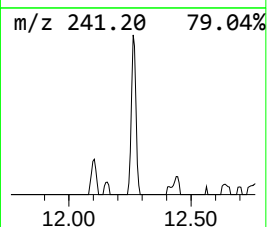
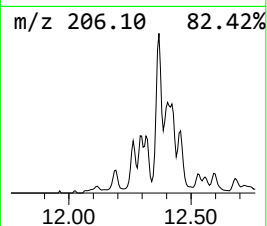
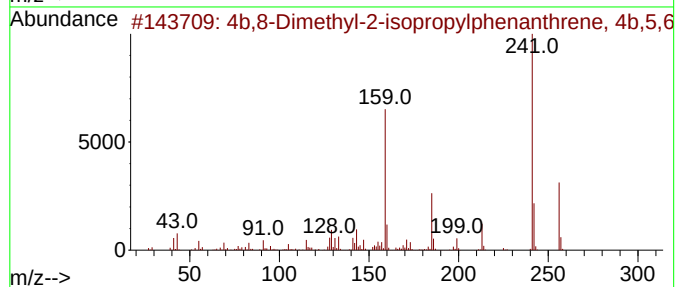
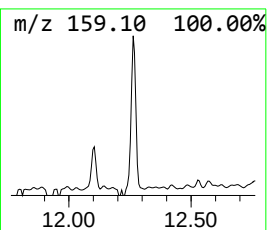
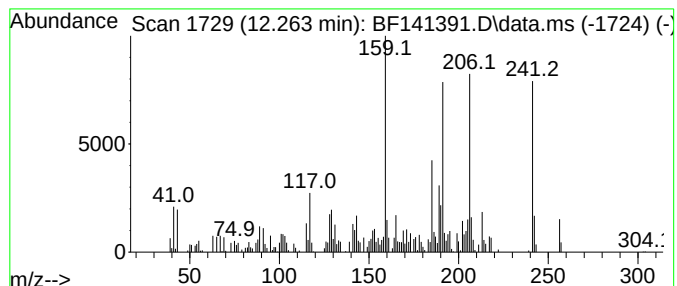
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4b,8-Dimethyl-2-isopropylph... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.263	2.86 ng	82065	Phenanthrene-d10	11.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	94
2			1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	72
3			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	62
4			Anthracene, 2-ethyl-	206	C16H14	052251-71-5	46
5			1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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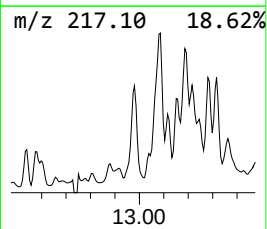
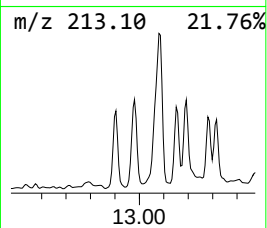
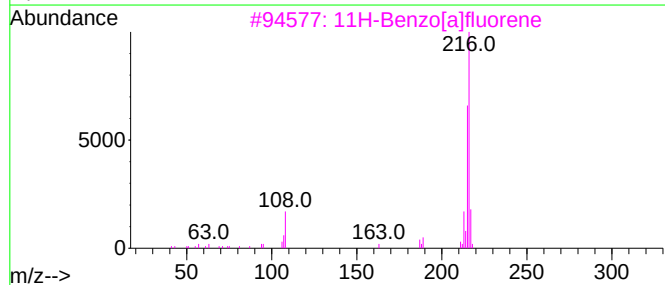
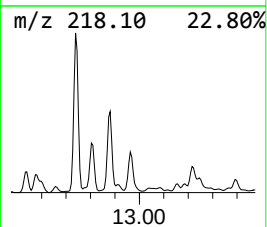
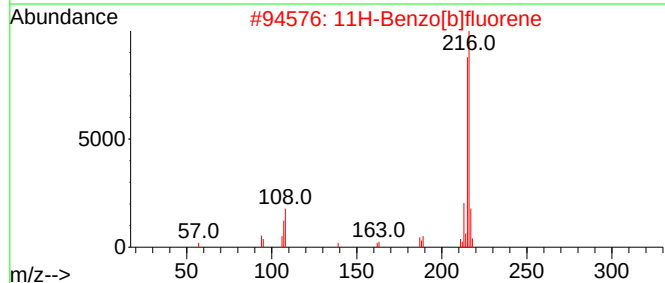
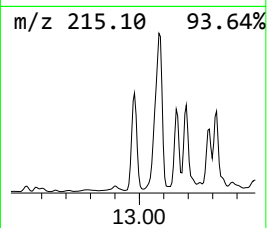
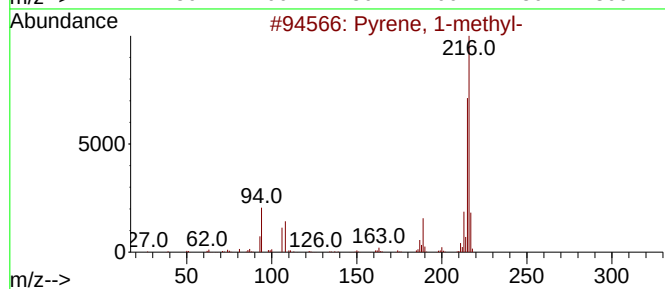
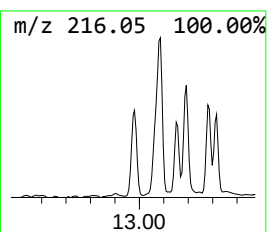
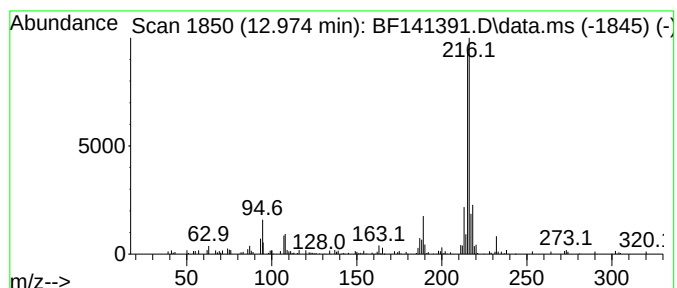
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.974	2.04 ng	108152	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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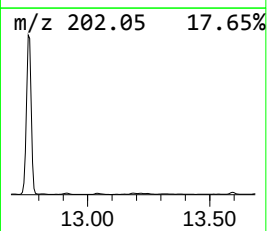
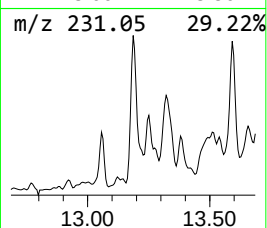
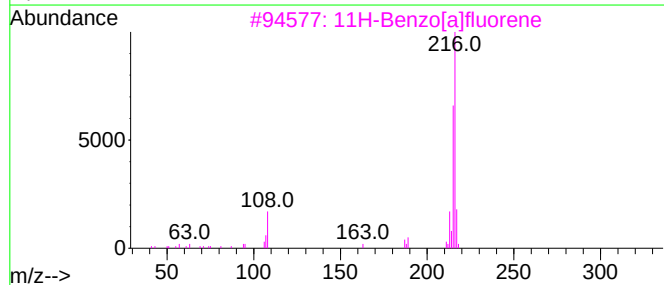
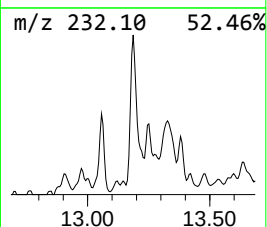
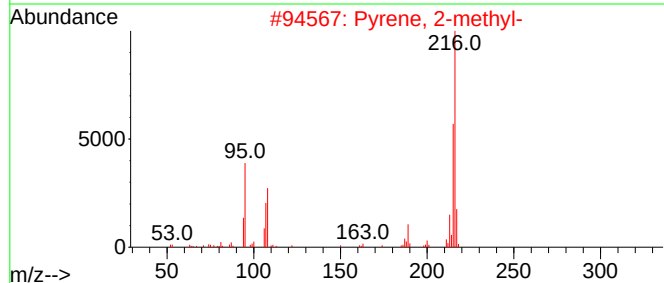
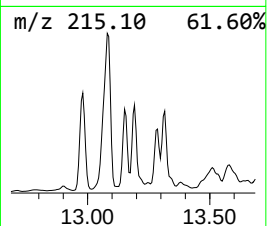
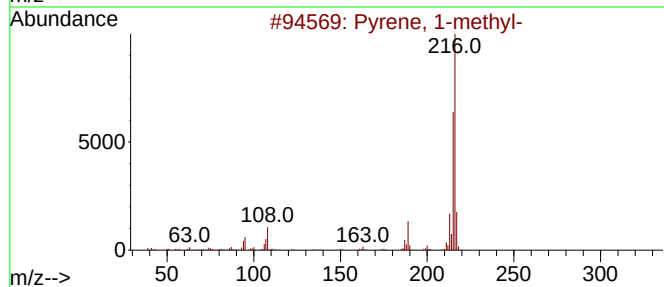
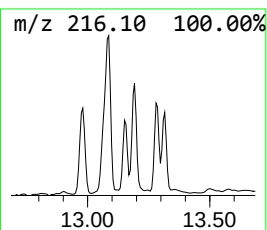
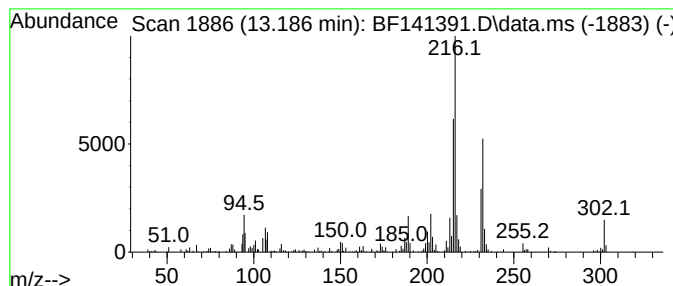
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.186	4.09 ng	216803	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	78
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	49
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	46
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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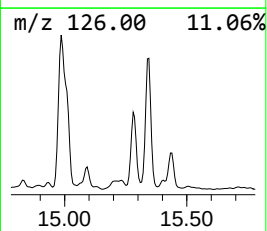
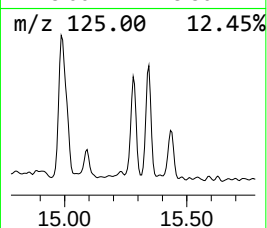
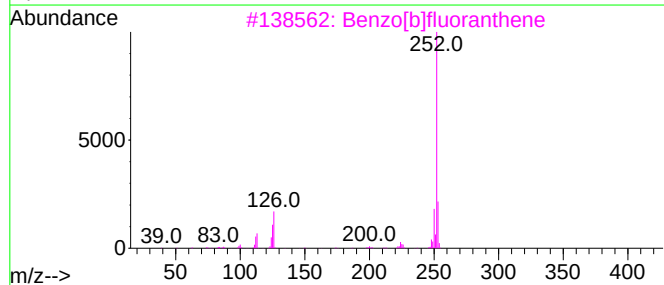
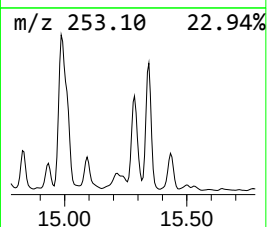
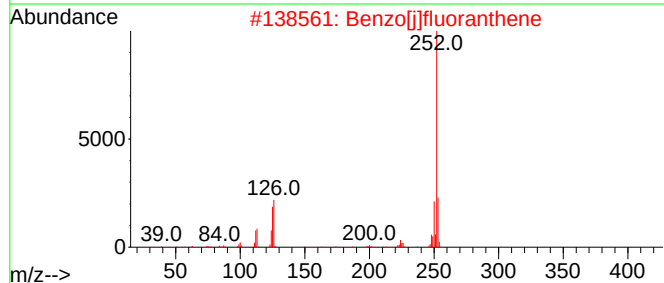
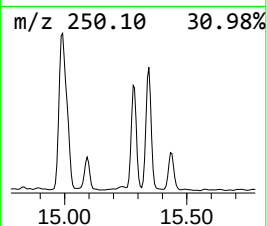
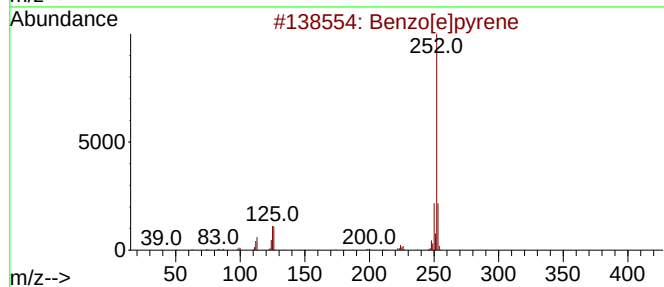
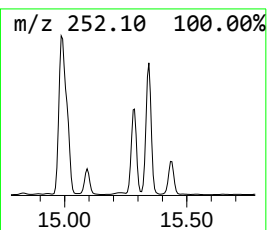
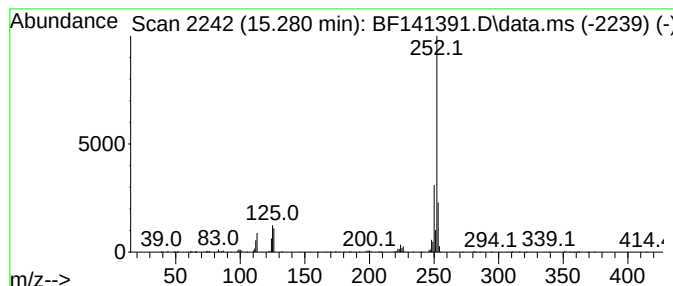
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.280	13.95 ng	236429	Perylene-d12	15.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
Data File : BF141391.D
Acq On : 04 Feb 2025 01:16
Operator : RC/JU
Sample : Q1216-19 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JPP-26.2-012825

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.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
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Anthracene, 9-m...	11.816	2.9	ng	81977	4	11.316	572937	20.0
Phenanthrene, 2...	11.845	6.0	ng	171738	4	11.316	572937	20.0
4H-Cyclopenta[d...	11.927	7.1	ng	203367	4	11.316	572937	20.0
Naphthalene, 2-...	12.116	2.7	ng	78035	4	11.316	572937	20.0
4b,8-Dimethyl-2...	12.263	2.9	ng	82065	4	11.316	572937	20.0
Pyrene, 1-methyl-	12.974	2.0	ng	108152	5	13.957	1059450	20.0
Pyrene, 2-methyl-	13.186	4.1	ng	216803	5	13.957	1059450	20.0
Benzo[e]pyrene	15.280	13.9	ng	236429	6	15.404	338983	20.0