

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041620\
 Data File : BF120026.D
 Acq On : 16 Apr 2020 10:46
 Operator : MA/SJ
 Sample : L2230-51 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-9-17

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF040820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.316	546	549	557	rVB	752268	694126	48.35%	2.912%
2	6.345	721	724	739	rBV	734145	711161	49.53%	2.984%
3	6.722	784	788	791	rBV	1181499	915028	63.73%	3.839%
4	6.875	811	814	818	rBV	799474	631601	43.99%	2.650%
5	7.281	879	883	889	rBV	611369	485679	33.83%	2.038%
6	8.010	1002	1007	1009	rBV	1326240	1202116	83.73%	5.044%
7	8.028	1009	1010	1014	rVB	93478	52984	3.69%	0.222%
8	8.722	1122	1128	1131	rBV2	52143	44948	3.13%	0.189%
9	8.822	1141	1145	1148	rBV	40887	31286	2.18%	0.131%
10	9.086	1186	1190	1193	rBV	1174759	943792	65.74%	3.960%
11	9.186	1200	1207	1209	rBV2	19333	32801	2.28%	0.138%
12	9.345	1227	1234	1239	rVB	23524	36093	2.51%	0.151%
13	9.422	1244	1247	1250	rBV	44684	43786	3.05%	0.184%
14	9.481	1254	1257	1262	rVB	138217	120086	8.36%	0.504%
15	9.539	1262	1267	1269	rBV3	27172	26913	1.87%	0.113%
16	9.622	1279	1281	1288	rVB	57975	44409	3.09%	0.186%
17	9.763	1299	1305	1308	rBV	1600541	1293809	90.12%	5.428%
18	9.792	1308	1310	1314	rVB	120612	111462	7.76%	0.468%
19	9.969	1337	1340	1344	rVV	98134	84981	5.92%	0.357%
20	10.081	1354	1359	1362	rBV2	32635	39018	2.72%	0.164%
21	10.175	1371	1375	1377	rBV2	31978	38645	2.69%	0.162%
22	10.310	1395	1398	1404	rBV	140699	126664	8.82%	0.531%
23	10.469	1422	1425	1430	rVB	47741	45824	3.19%	0.192%
24	10.551	1435	1439	1443	rVV	567802	519603	36.19%	2.180%
25	10.675	1458	1460	1467	rVB2	45451	62804	4.37%	0.264%
26	10.863	1485	1492	1494	rBV4	43919	75637	5.27%	0.317%
27	10.886	1494	1496	1499	rVV2	38480	38271	2.67%	0.161%
28	10.939	1503	1505	1508	rVV3	27283	29358	2.04%	0.123%
29	10.986	1508	1513	1515	rVV3	36541	51690	3.60%	0.217%
30	11.010	1515	1517	1522	rVV2	38948	51338	3.58%	0.215%
31	11.057	1522	1525	1529	rVV2	53443	56662	3.95%	0.238%
32	11.098	1529	1532	1535	rVV2	40681	59948	4.18%	0.252%
33	11.145	1538	1540	1546	rVB	84944	88691	6.18%	0.372%
34	11.251	1554	1558	1560	rBV	1353944	1237373	86.19%	5.192%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.275	1560	1562	1565	rVV	1263540	909864	63.37%	3.817%
36	11.322	1565	1570	1574	rVB	290960	258568	18.01%	1.085%
37	11.357	1574	1576	1580	rBV2	53543	60300	4.20%	0.253%
38	11.480	1592	1597	1599	rBV2	93018	98165	6.84%	0.412%
39	11.504	1599	1601	1605	rVV4	30969	38737	2.70%	0.163%
40	11.551	1606	1609	1612	rVV3	52642	55169	3.84%	0.231%
41	11.580	1612	1614	1619	rVB3	48330	45631	3.18%	0.191%
42	11.751	1639	1643	1645	rVV	192428	164041	11.43%	0.688%
43	11.780	1645	1648	1653	rVV2	407590	424338	29.56%	1.780%
44	11.827	1653	1656	1659	rVV	99468	102007	7.11%	0.428%
45	11.863	1659	1662	1664	rVV	290807	287361	20.02%	1.206%
46	11.886	1664	1666	1671	rVB	131516	115288	8.03%	0.484%
47	11.986	1681	1683	1689	rBV5	19605	38121	2.66%	0.160%
48	12.045	1690	1693	1695	rVV	103566	96010	6.69%	0.403%
49	12.069	1695	1697	1702	rVB2	72161	63495	4.42%	0.266%
50	12.198	1716	1719	1722	rVB2	70303	65627	4.57%	0.275%
51	12.227	1722	1724	1726	rBV	43157	40732	2.84%	0.171%
52	12.251	1726	1728	1731	rVB2	47971	37995	2.65%	0.159%
53	12.304	1734	1737	1740	rVV	107785	113325	7.89%	0.475%
54	12.339	1740	1743	1748	rVB3	71991	127991	8.91%	0.537%
55	12.386	1748	1751	1756	rVB2	82935	95716	6.67%	0.402%
56	12.463	1760	1764	1767	rBV	1487050	1270363	88.48%	5.330%
57	12.510	1767	1772	1778	rVB6	59180	92355	6.43%	0.387%
58	12.569	1778	1782	1788	rBV4	79683	139557	9.72%	0.586%
59	12.692	1797	1803	1806	rBV	1174221	1193985	83.16%	5.010%
60	12.739	1809	1811	1817	rVB2	60785	83258	5.80%	0.349%
61	12.810	1820	1823	1825	rBV	90518	92446	6.44%	0.388%
62	12.833	1825	1827	1834	rVB	817801	774910	53.97%	3.251%
63	12.910	1835	1840	1843	rBV2	90383	144118	10.04%	0.605%
64	12.998	1852	1855	1856	rBV3	50817	58217	4.05%	0.244%
65	13.022	1856	1859	1862	rVB2	178196	173241	12.07%	0.727%
66	13.086	1866	1870	1872	rBV2	152529	133196	9.28%	0.559%
67	13.121	1872	1876	1879	rBV	133953	140987	9.82%	0.592%
68	13.216	1889	1892	1895	rVB	78242	65111	4.54%	0.273%
69	13.245	1895	1897	1901	rBV2	79369	72831	5.07%	0.306%
70	13.304	1905	1907	1910	rVV3	38313	30437	2.12%	0.128%
71	13.527	1942	1945	1950	rVB	72661	91665	6.38%	0.385%

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72	13.633	1960	1963	1966	rBV3	82709	100720	7.02%	0.423%
73	13.669	1966	1969	1971	rVV	82203	79512	5.54%	0.334%
74	13.686	1971	1972	1978	rVB2	59424	69524	4.84%	0.292%
75	13.733	1978	1980	1982	rBV	53440	44861	3.12%	0.188%
76	13.886	1999	2006	2009	rBV2	1037665	1435701	100.00%	6.024%
77	13.916	2009	2011	2013	rVB	481528	361973	25.21%	1.519%
78	13.992	2022	2024	2027	rVB	77827	64061	4.46%	0.269%
79	14.069	2035	2037	2041	rVB2	36368	39691	2.76%	0.167%
80	14.210	2059	2061	2064	rBV3	26448	27051	1.88%	0.113%
81	14.239	2064	2066	2068	rVV	40834	38546	2.68%	0.162%
82	14.274	2068	2072	2075	rVV	117374	141643	9.87%	0.594%
83	14.310	2075	2078	2081	rVB	66625	66136	4.61%	0.277%
84	14.368	2086	2088	2091	rVV2	95381	96639	6.73%	0.405%
85	14.398	2091	2093	2095	rVV	62747	64648	4.50%	0.271%
86	14.421	2095	2097	2101	rVB3	65165	65549	4.57%	0.275%
87	14.910	2175	2180	2183	rBV	390340	633116	44.10%	2.656%
88	14.992	2190	2194	2195	rBV2	69976	87765	6.11%	0.368%
89	15.192	2224	2228	2234	rVB	235242	307366	21.41%	1.290%
90	15.251	2234	2238	2244	rBV	356459	427626	29.79%	1.794%
91	15.315	2244	2249	2252	rVV	593834	711223	49.54%	2.984%
92	15.345	2252	2254	2260	rVB3	129794	200907	13.99%	0.843%
93	15.768	2322	2326	2331	rBV	64467	90398	6.30%	0.379%
94	15.921	2346	2352	2354	rBV7	27812	63526	4.42%	0.267%
95	16.245	2403	2407	2411	rBV4	29464	51879	3.61%	0.218%
96	16.692	2478	2483	2491	rVB2	181237	355955	24.79%	1.493%
97	16.862	2510	2512	2517	rVB2	40771	57619	4.01%	0.242%
98	16.921	2518	2522	2526	rBV4	23120	47629	3.32%	0.200%
99	17.109	2549	2554	2564	rBV	127209	250758	17.47%	1.052%
100	17.327	2588	2591	2592	rBV3	26355	28636	1.99%	0.120%

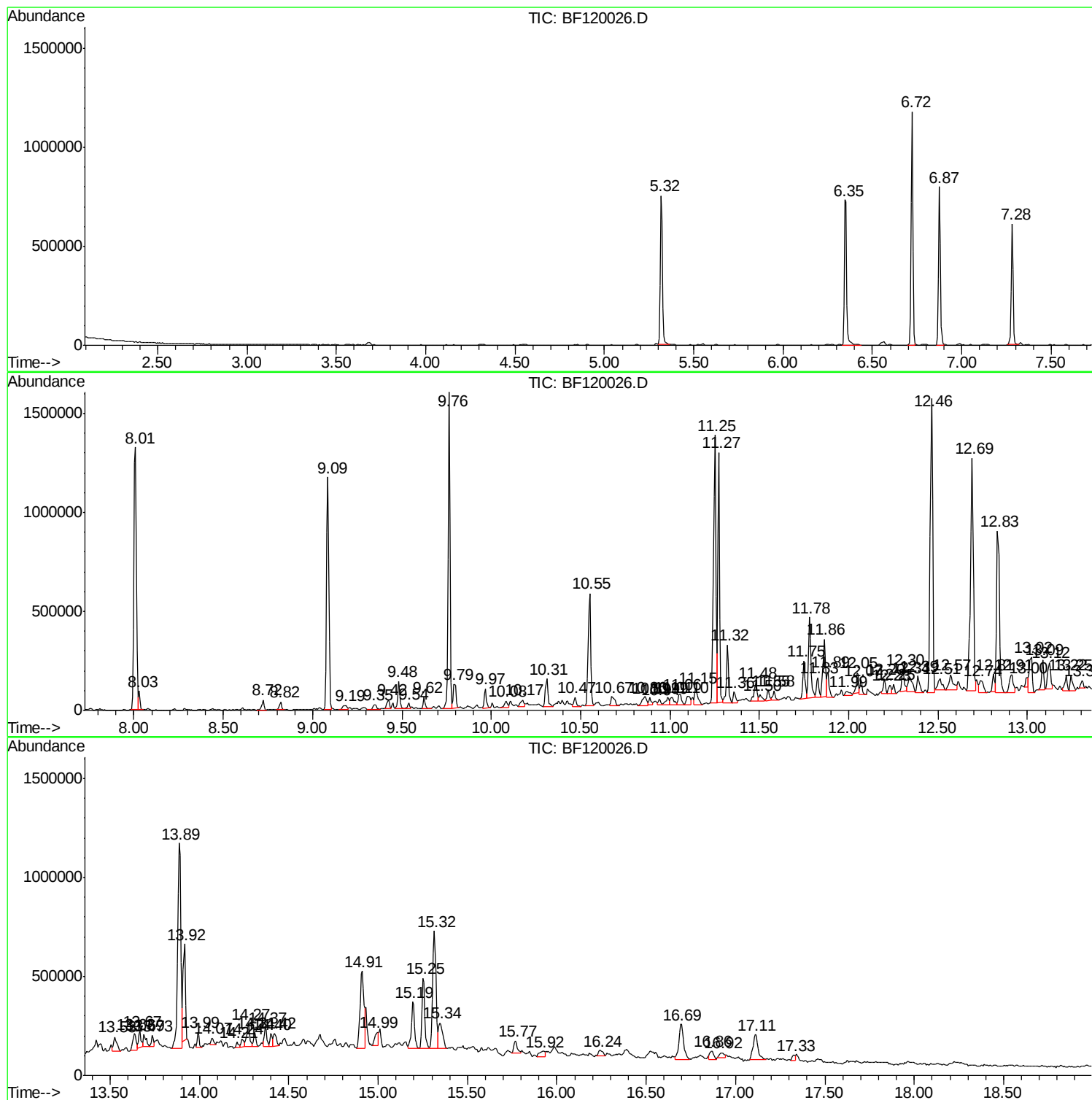
Sum of corrected areas: 23834403

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Instrument :
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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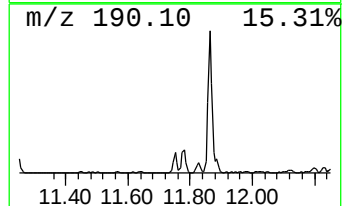
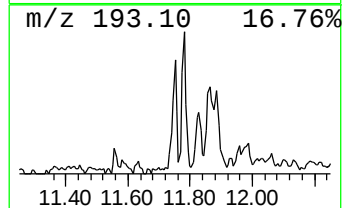
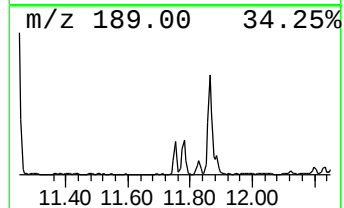
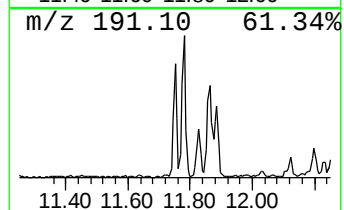
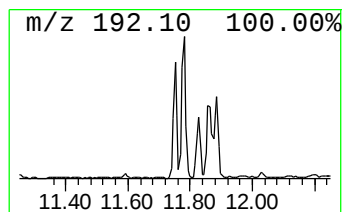
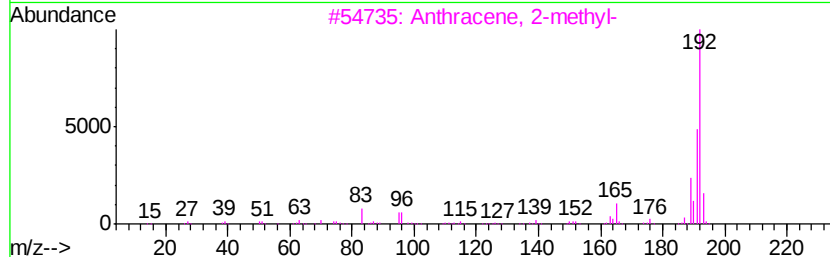
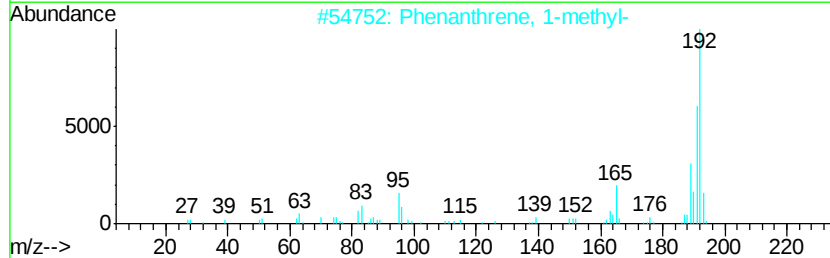
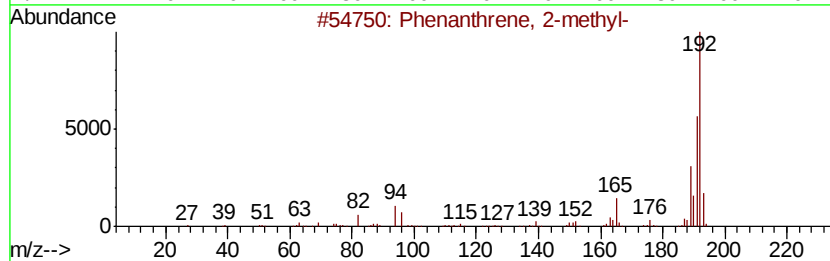
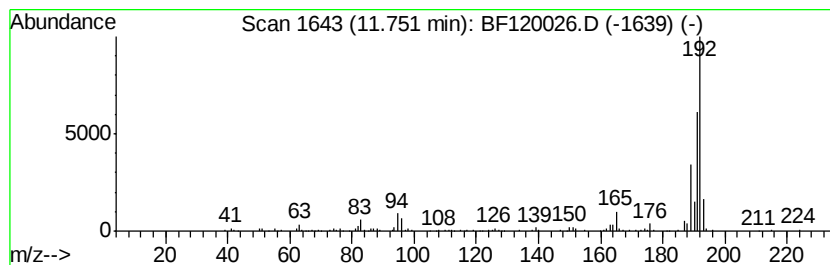
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	2.65 ng	164041	Phenanthrene-d10	11.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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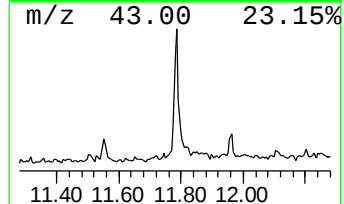
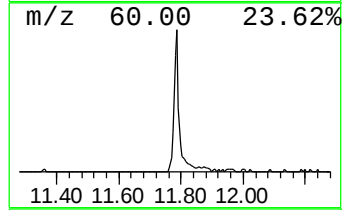
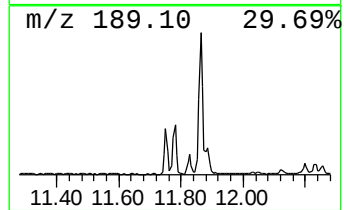
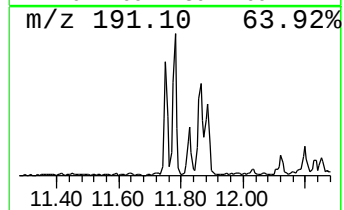
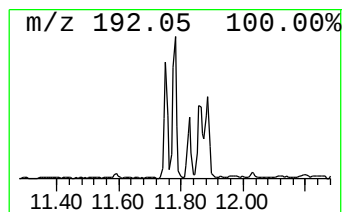
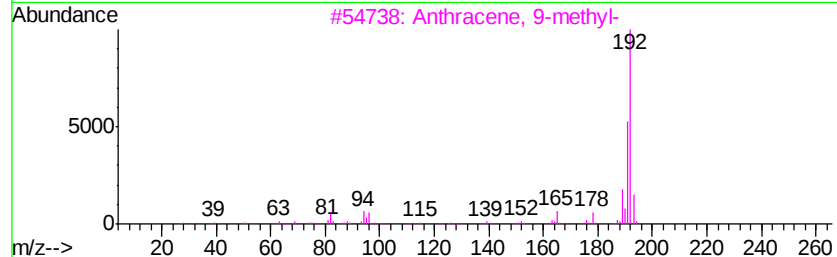
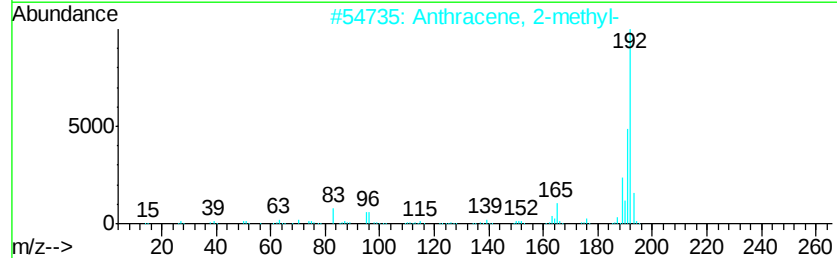
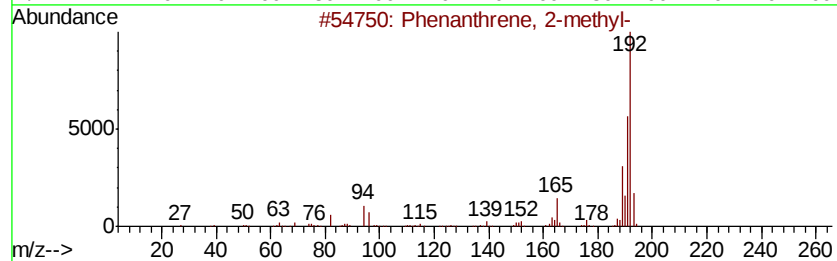
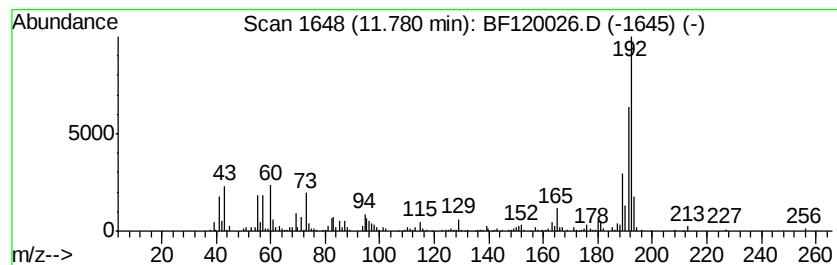
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.78	6.86 ng	424338	Phenanthrene-d10	11.25	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	92
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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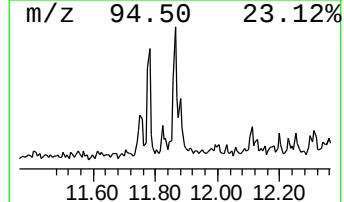
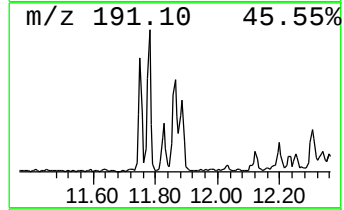
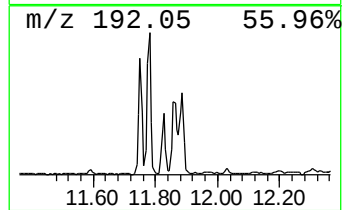
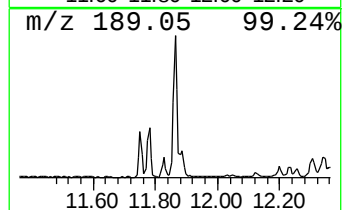
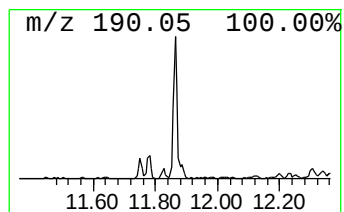
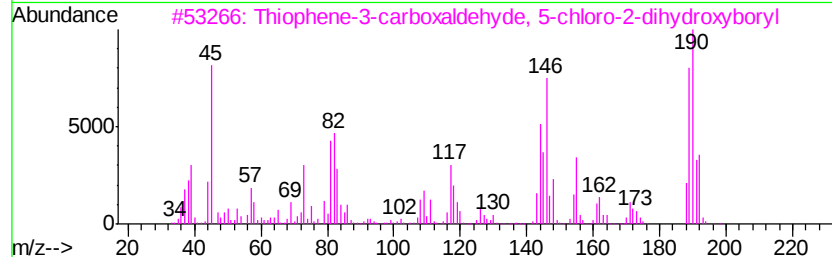
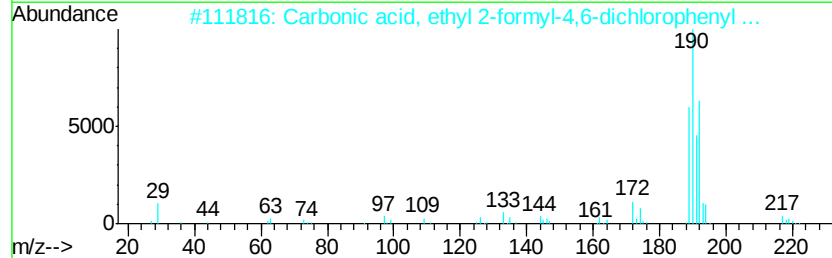
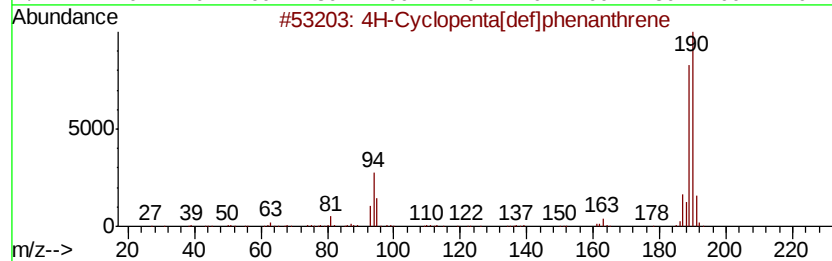
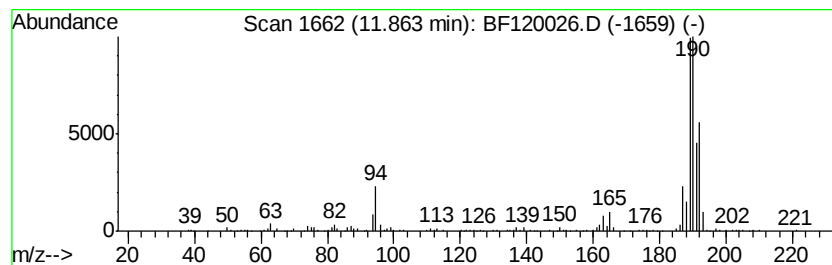
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.86	4.64 ng	287361	Phenanthrene-d10	11.25	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
4	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120026.D
Acq On : 16 Apr 2020 10:46
Operator : MA/SJ
Sample : L2230-51 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

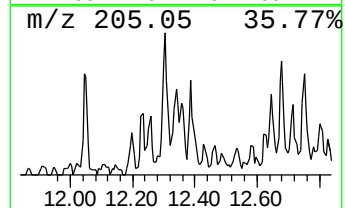
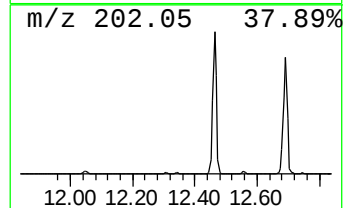
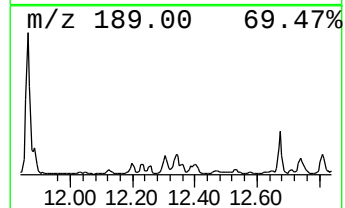
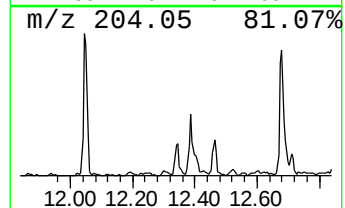
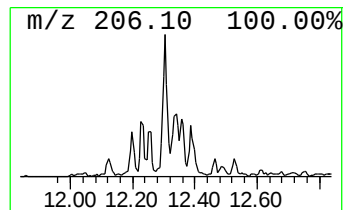
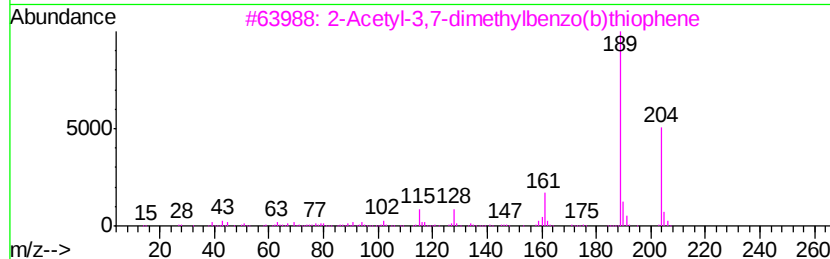
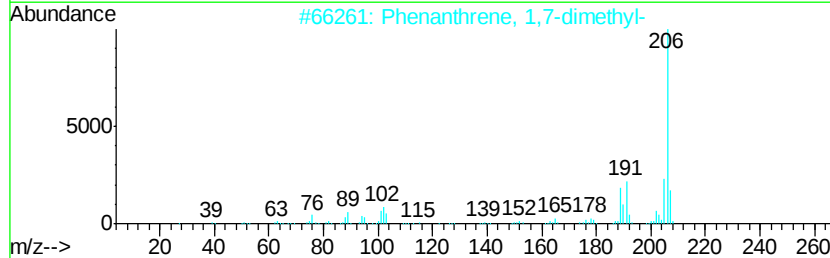
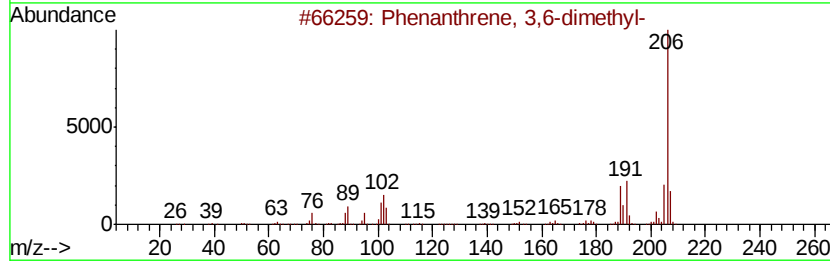
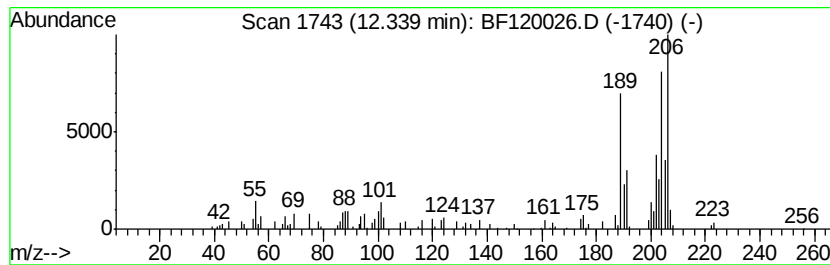
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown12.34 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.34	2.07 ng	127991	Phenanthrene-d10	11.25	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	41
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	38
3	2-Acetyl-3,7-dimethylbenzo(b)thi...	204	C12H12OS	006179-06-2	35
4	4-Pyridinol 3,5-dichloro-2-ethyl...	205	C8H9Cl2NO	1000253-55-0	35
5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120026.D
Acq On : 16 Apr 2020 10:46
Operator : MA/SJ
Sample : L2230-51 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

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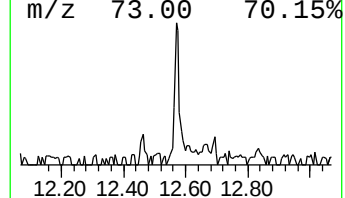
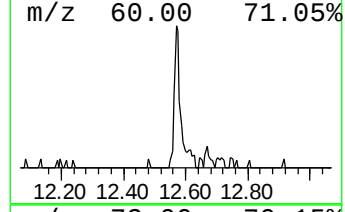
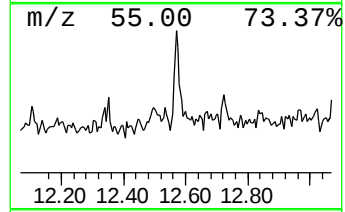
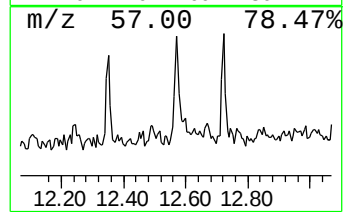
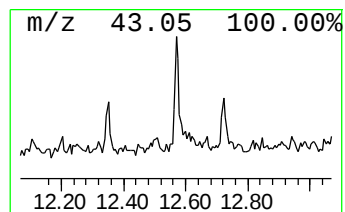
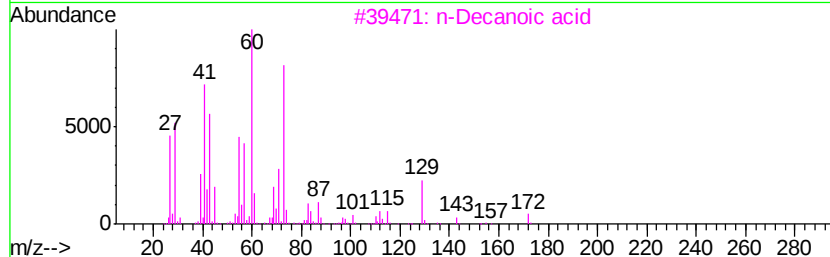
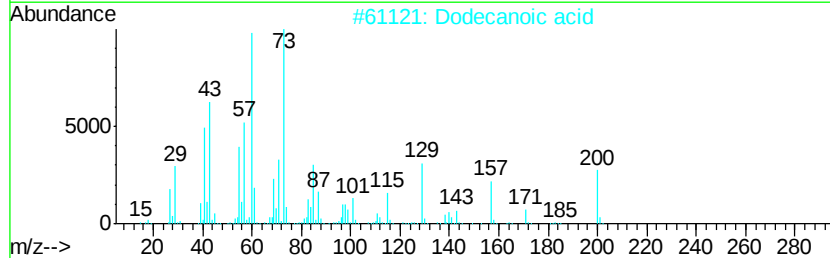
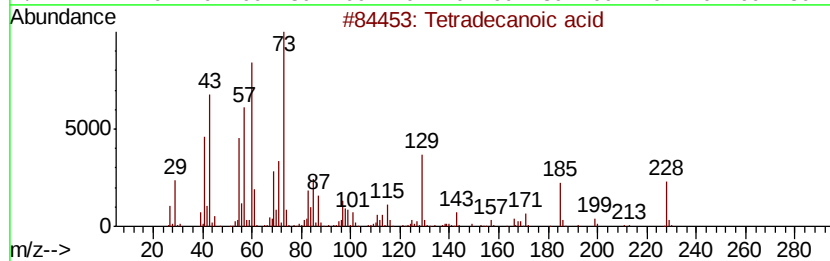
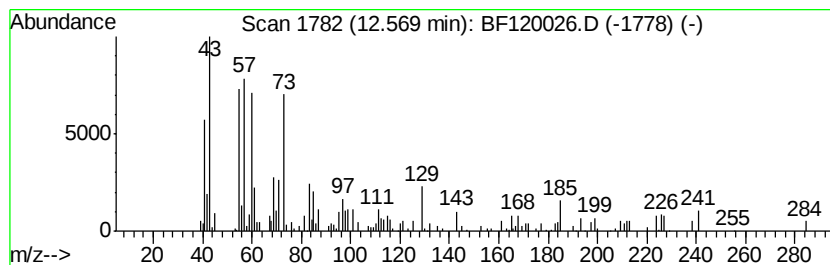
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetradecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	2.26 ng	139557	Phenanthrene-d10	11.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
2		Dodecanoic acid	200	C12H24O2	000143-07-7	50
3		n-Decanoic acid	172	C10H20O2	000334-48-5	50
4		Octadecanoic acid	284	C18H36O2	000057-11-4	46
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120026.D
Acq On : 16 Apr 2020 10:46
Operator : MA/SJ
Sample : L2230-51 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

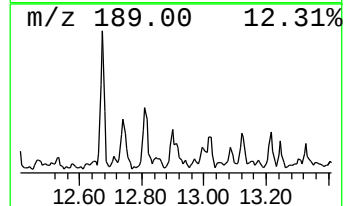
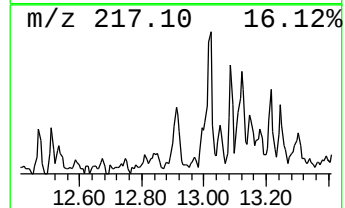
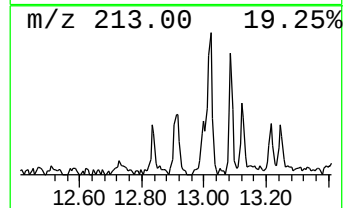
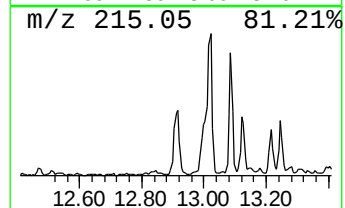
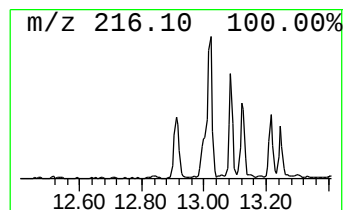
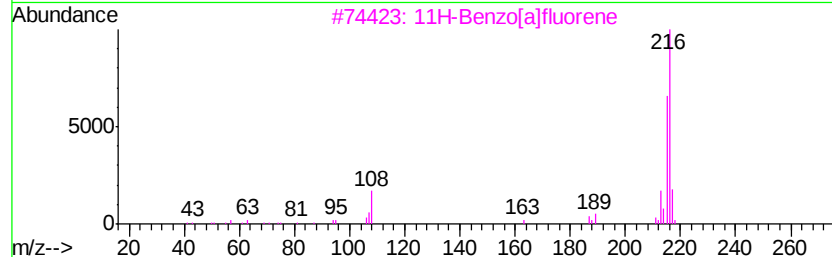
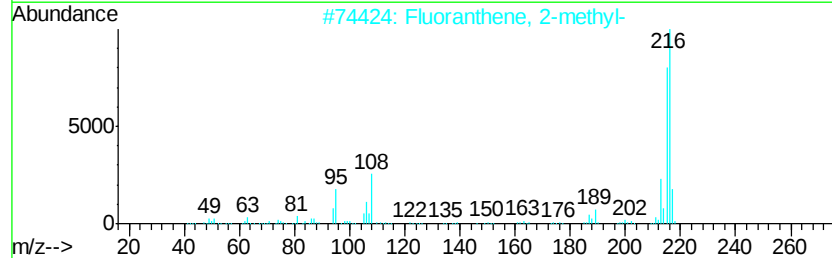
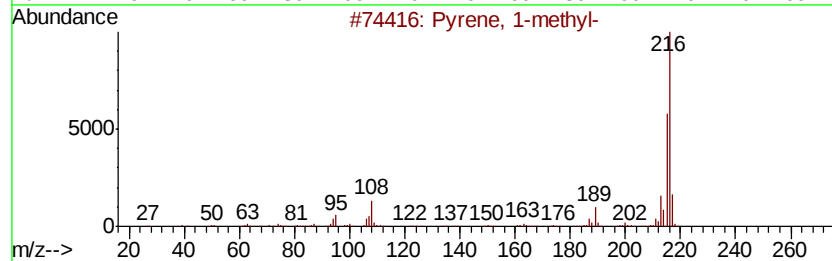
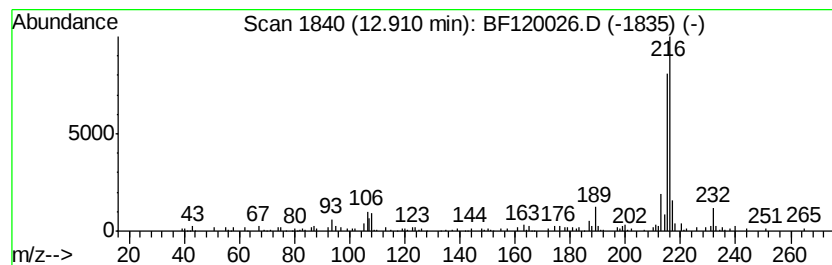
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.91	2.01 ng	144118	Chrysene-d12	13.89		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	89
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120026.D
Acq On : 16 Apr 2020 10:46
Operator : MA/SJ
Sample : L2230-51 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

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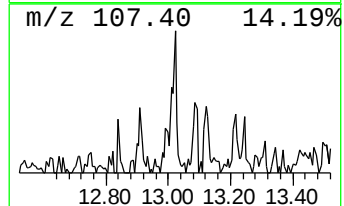
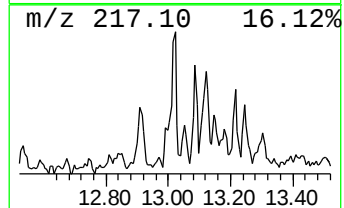
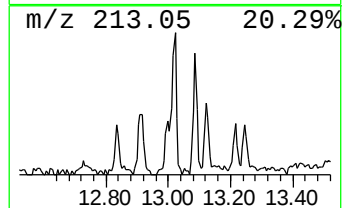
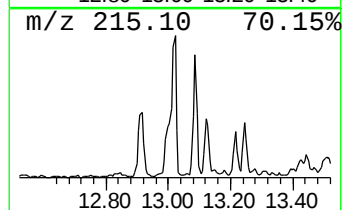
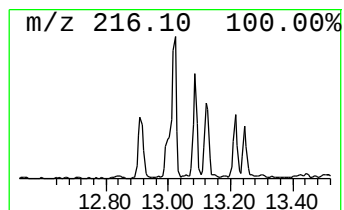
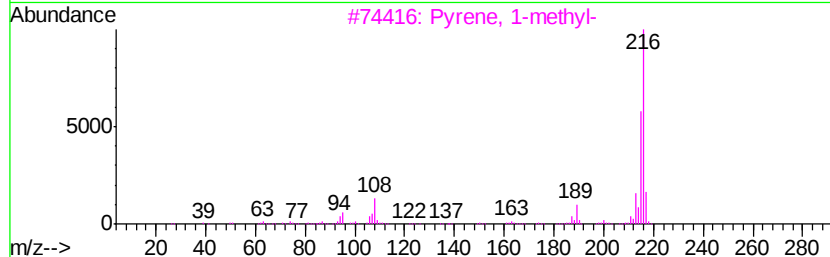
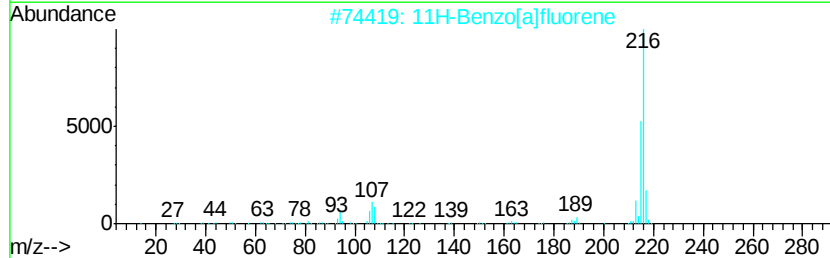
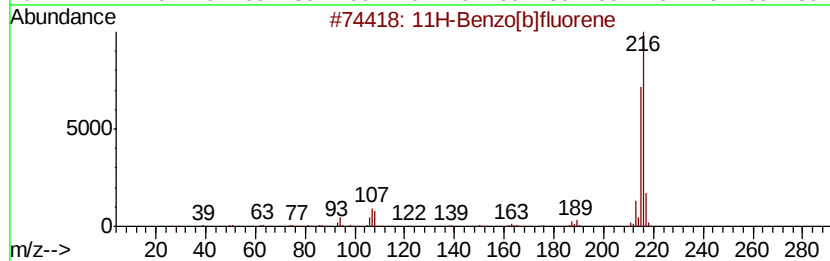
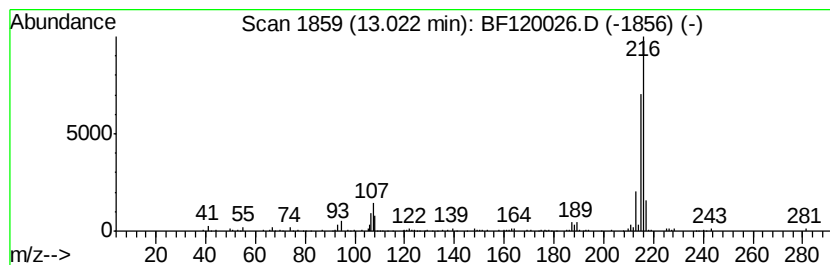
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	2.41 ng	173241	Chrysene-d12	13.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	80
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120026.D
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Sample : L2230-51 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

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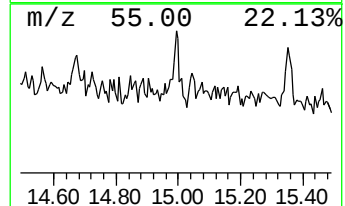
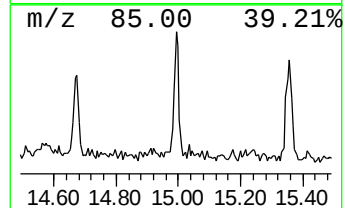
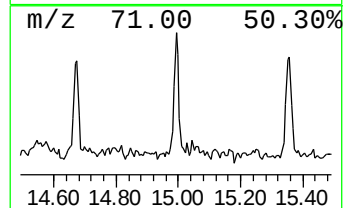
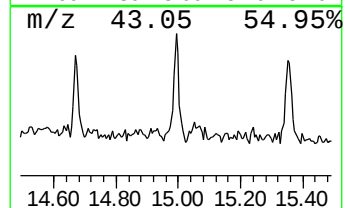
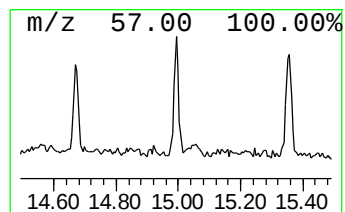
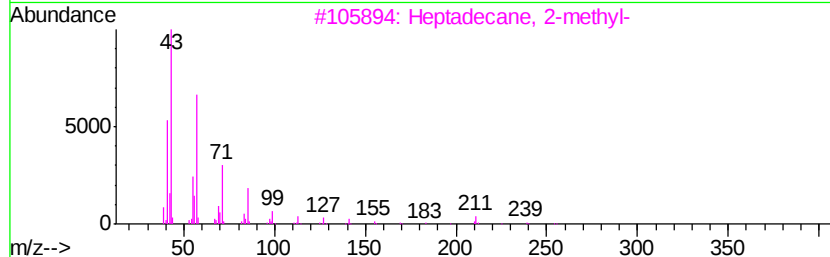
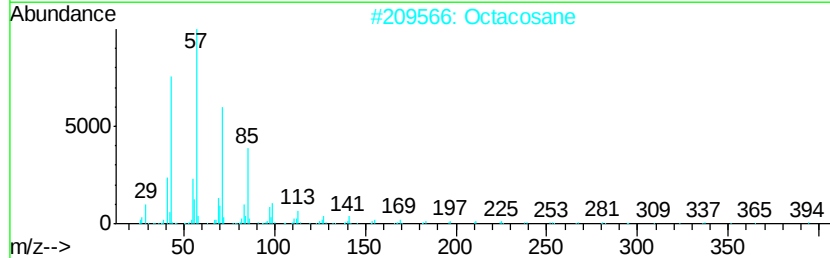
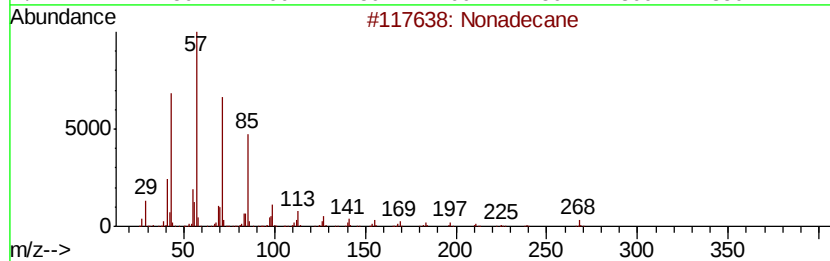
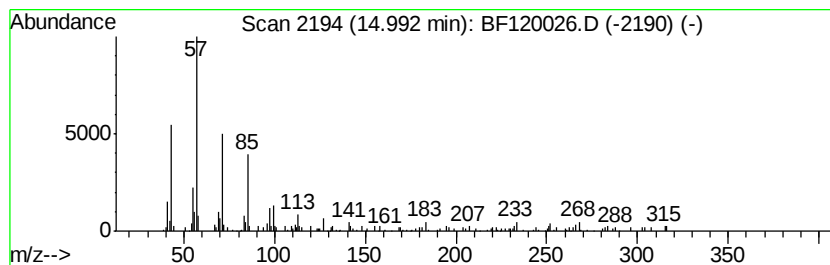
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	2.47 ng	87765	Perylene-d12	15.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	90
2	Octacosane		394	C28H58	000630-02-4	87
3	Heptadecane, 2-methyl-		254	C18H38	001560-89-0	81
4	Docosane		310	C22H46	000629-97-0	74
5	Tetracosane		338	C24H50	000646-31-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
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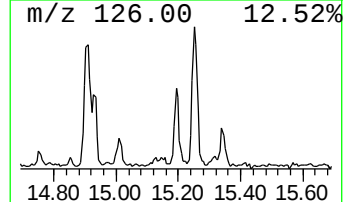
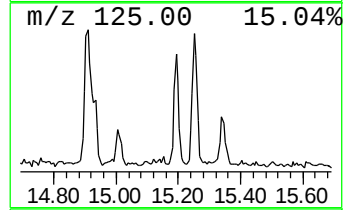
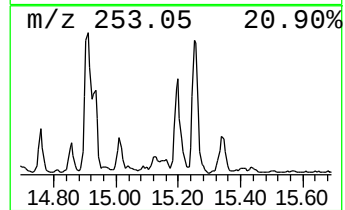
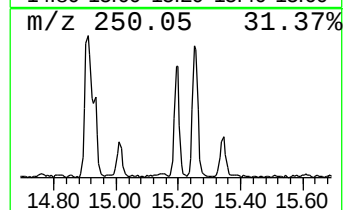
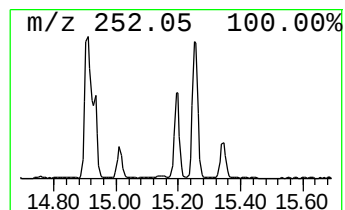
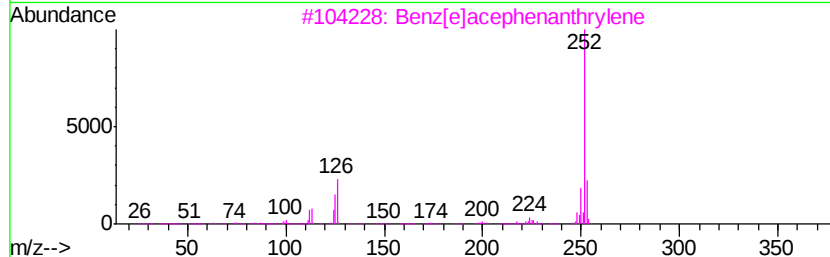
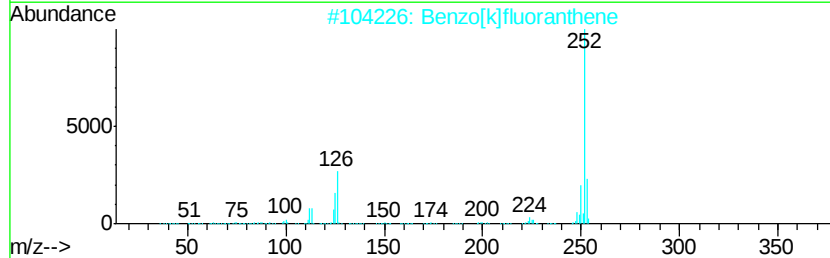
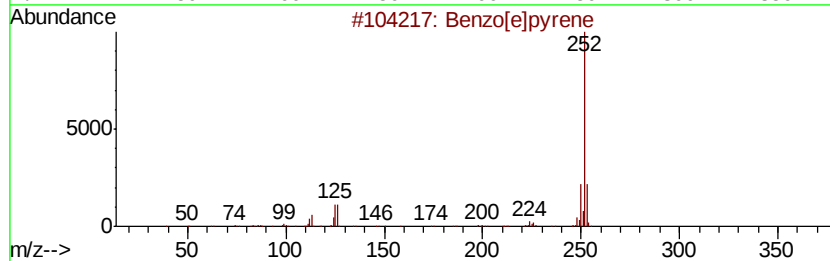
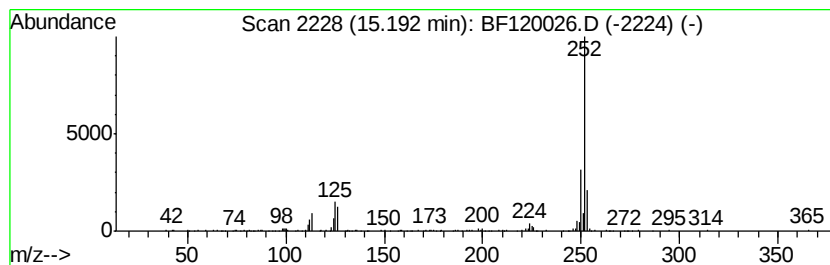
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.19	8.64 ng	307366	Perylene-d12	15.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	94
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3	Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
4	Benzo[a]pyrene	252	C20H12	000050-32-8	93
5	Perylene	252	C20H12	000198-55-0	91



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ALS Vial : 34 Sample Multiplier: 1

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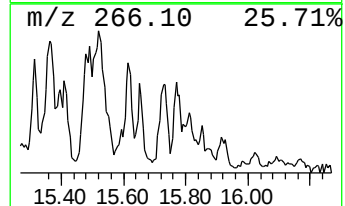
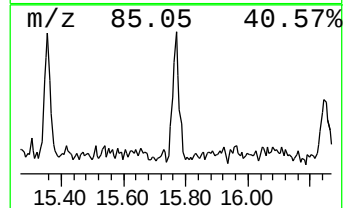
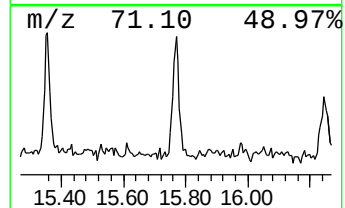
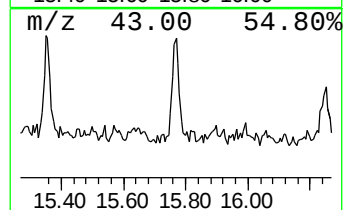
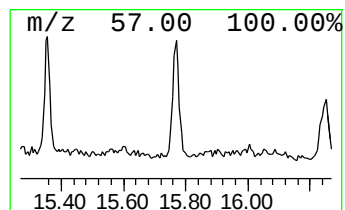
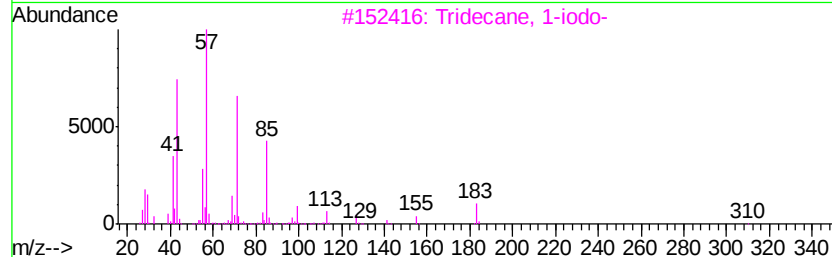
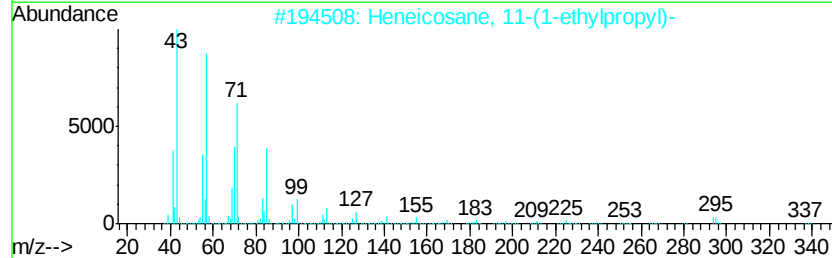
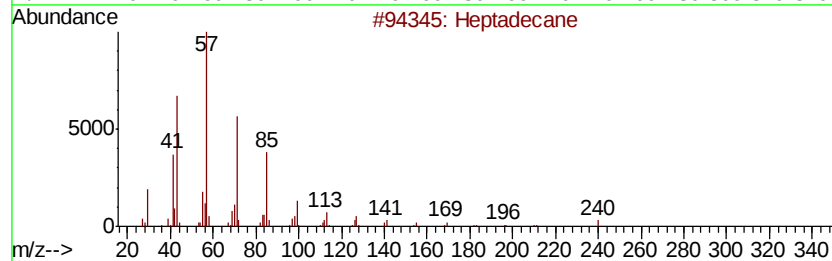
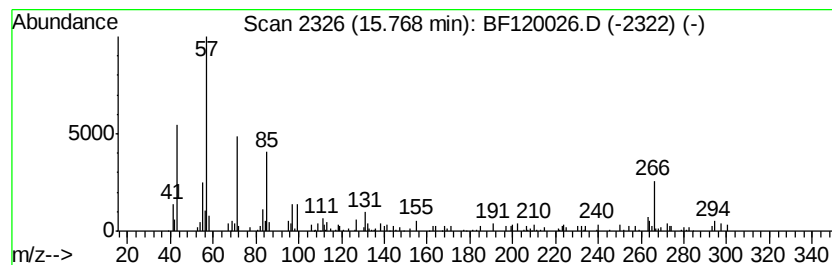
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	2.54 ng	90398	Perylene-d12	15.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	60
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	49
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	49
4		Hexadecane	226	C16H34	000544-76-3	49
5		Pentacosane	352	C25H52	000629-99-2	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041620\
 Data File : BF120026.D
 Acq On : 16 Apr 2020 10:46
 Operator : MA/SJ
 Sample : L2230-51 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-9-17

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 2-m...	11.75	2.6	ng	164041	4	11.25	1237370	20.0
Anthracene, 2-met...	11.78	6.9	ng	424338	4	11.25	1237370	20.0
4H-Cyclopenta[def...	11.86	4.6	ng	287361	4	11.25	1237370	20.0
unknown12.34	12.34	2.1	ng	127991	4	11.25	1237370	20.0
Tetradecanoic acid	12.57	2.3	ng	139557	4	11.25	1237370	20.0
Pyrene, 1-methyl-	12.91	2.0	ng	144118	5	13.89	1435700	20.0
11H-Benzo[b]fluorene	13.02	2.4	ng	173241	5	13.89	1435700	20.0
Nonadecane	14.99	2.5	ng	87765	6	15.32	711223	20.0
Benzo[e]pyrene	15.19	8.6	ng	307366	6	15.32	711223	20.0
Heptadecane	15.77	2.5	ng	90398	6	15.32	711223	20.0