

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102919\
 Data File : BF117587.D
 Acq On : 29 Oct 2019 13:37
 Operator : JU
 Sample : K5503-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 172-80-(0-2)

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-BF101819.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	5.357	553	556	572	rBV	446145	631681	84.26%	7.108%
2	6.386	727	731	748	rBV	539005	732008	97.65%	8.236%
3	6.504	748	751	766	rVB	755769	749651	100.00%	8.435%
4	6.728	786	789	799	rBV	295419	265020	35.35%	2.982%
5	6.881	812	815	829	rBV	536332	511849	68.28%	5.759%
6	7.292	882	885	909	rBV	326530	419276	55.93%	4.718%
7	8.010	1003	1007	1031	rBV	289256	330364	44.07%	3.717%
8	9.080	1186	1189	1201	rBV	645727	601092	80.18%	6.763%
9	9.480	1254	1257	1265	rBV	70242	67958	9.07%	0.765%
10	9.763	1301	1305	1309	rBV	311676	307422	41.01%	3.459%
11	9.792	1309	1310	1319	rVB	17074	14670	1.96%	0.165%
12	10.316	1395	1399	1404	rBB	13551	14515	1.94%	0.163%
13	10.557	1436	1440	1452	rBV	261078	289250	38.58%	3.255%
14	11.145	1537	1540	1545	rVB2	7067	7726	1.03%	0.087%
15	11.245	1554	1557	1560	rBV	260314	272472	36.35%	3.066%
16	11.269	1560	1561	1567	rVV	184942	174467	23.27%	1.963%
17	11.327	1567	1571	1579	rVB	34945	40901	5.46%	0.460%
18	11.498	1595	1600	1604	rBV3	7513	11007	1.47%	0.124%
19	11.751	1640	1643	1645	rBV	32586	28345	3.78%	0.319%
20	11.780	1645	1648	1654	rVV3	68712	83324	11.12%	0.938%
21	11.827	1654	1656	1659	rVV	14631	16008	2.14%	0.180%
22	11.863	1659	1662	1664	rVV2	57277	57520	7.67%	0.647%
23	11.886	1664	1666	1673	rVB	31788	28605	3.82%	0.322%
24	12.045	1690	1693	1699	rBV	20112	19520	2.60%	0.220%
25	12.092	1699	1701	1704	rBV	10940	10002	1.33%	0.113%
26	12.298	1733	1736	1739	rBV	22423	23226	3.10%	0.261%
27	12.333	1739	1742	1748	rVB4	13397	21781	2.91%	0.245%
28	12.398	1748	1753	1759	rVB2	14503	19737	2.63%	0.222%
29	12.463	1761	1764	1768	rBV	315436	279510	37.29%	3.145%
30	12.527	1773	1775	1779	rVV4	8987	12256	1.63%	0.138%
31	12.563	1779	1781	1786	rVV5	8712	12178	1.62%	0.137%
32	12.615	1787	1790	1793	rVV	16537	17478	2.33%	0.197%
33	12.692	1797	1803	1810	rVV	295563	326182	43.51%	3.670%
34	12.745	1810	1812	1817	rVB2	15247	17906	2.39%	0.201%

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35	12.827	1822	1826	1831	rBV	498986	423040	56.43%	4.760%
36	12.921	1836	1842	1845	rBV3	21136	35795	4.77%	0.403%
37	13.004	1852	1856	1857	rBV3	17803	21888	2.92%	0.246%
38	13.021	1857	1859	1862	rVB2	37566	33152	4.42%	0.373%
39	13.086	1867	1870	1873	rVV	28717	26406	3.52%	0.297%
40	13.127	1873	1877	1880	rVV2	32409	37164	4.96%	0.418%
41	13.180	1884	1886	1889	rVB3	8333	9100	1.21%	0.102%
42	13.221	1890	1893	1896	rBV	23181	23647	3.15%	0.266%
43	13.251	1896	1898	1903	rVV	24511	24021	3.20%	0.270%
44	13.398	1919	1923	1925	rBV4	17035	16840	2.25%	0.189%
45	13.445	1929	1931	1934	rVV3	9461	10098	1.35%	0.114%
46	13.510	1938	1942	1944	rBV5	12416	15498	2.07%	0.174%
47	13.545	1944	1948	1950	rVV	19115	17364	2.32%	0.195%
48	13.645	1961	1965	1967	rBV2	24614	30660	4.09%	0.345%
49	13.668	1967	1969	1971	rVB	20085	15477	2.06%	0.174%
50	13.698	1971	1974	1976	rBV	16720	14457	1.93%	0.163%
51	13.727	1976	1979	1982	rBV5	39800	56002	7.47%	0.630%
52	13.892	2002	2007	2010	rVV2	352494	466360	62.21%	5.247%
53	13.921	2010	2012	2019	rVB	152645	141384	18.86%	1.591%
54	13.998	2023	2025	2028	rVB	19711	18781	2.51%	0.211%
55	14.039	2029	2032	2034	rBV	24461	24099	3.21%	0.271%
56	14.162	2051	2053	2057	rVB5	7528	10514	1.40%	0.118%
57	14.286	2071	2074	2076	rVB	29710	28281	3.77%	0.318%
58	14.345	2081	2084	2088	rBV2	36914	43006	5.74%	0.484%
59	14.404	2092	2094	2097	rVB4	22378	23404	3.12%	0.263%
60	14.639	2132	2134	2138	rVB	48498	39419	5.26%	0.444%
61	14.709	2144	2146	2148	rVB3	14666	12333	1.65%	0.139%
62	14.921	2179	2182	2185	rBV	108453	149643	19.96%	1.684%
63	14.951	2185	2187	2192	rVB2	78737	99681	13.30%	1.122%
64	15.027	2197	2200	2203	rVB	27212	29604	3.95%	0.333%
65	15.209	2228	2231	2235	rBV	58652	63709	8.50%	0.717%
66	15.274	2238	2242	2245	rBV	75633	85607	11.42%	0.963%
67	15.327	2245	2251	2255	rVV	165466	232312	30.99%	2.614%
68	15.715	2314	2317	2321	rVB2	34845	42726	5.70%	0.481%
69	16.180	2393	2396	2401	rVB4	18307	28020	3.74%	0.315%
70	16.345	2422	2424	2430	rVB8	9534	17046	2.27%	0.192%
71	16.745	2488	2492	2500	rVB3	29523	62921	8.39%	0.708%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	17.180	2561	2566	2574	rVB3	19298	43122	5.75%	0.485%
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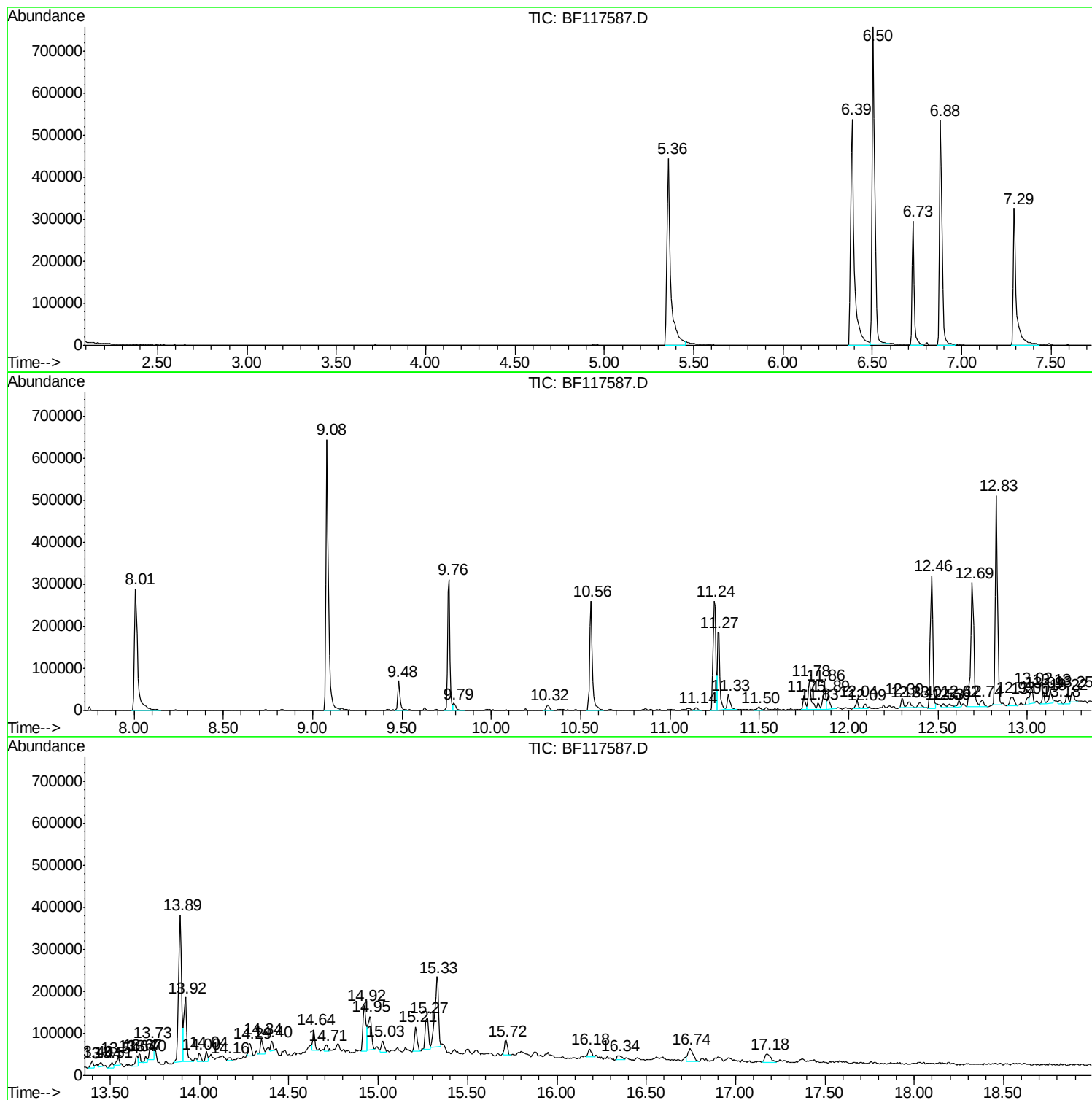
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Instrument :
BNA_F
ClientSampled :
172-80-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.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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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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172-80-(0-2)

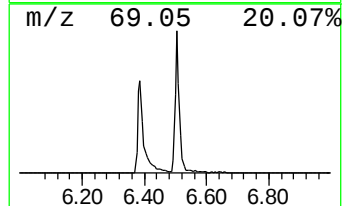
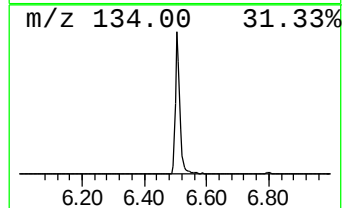
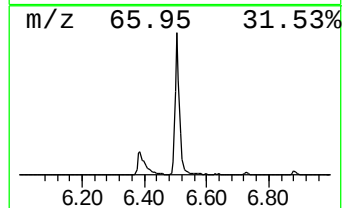
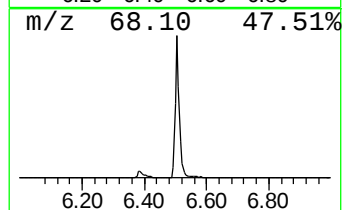
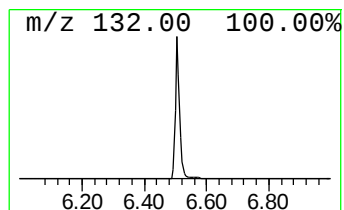
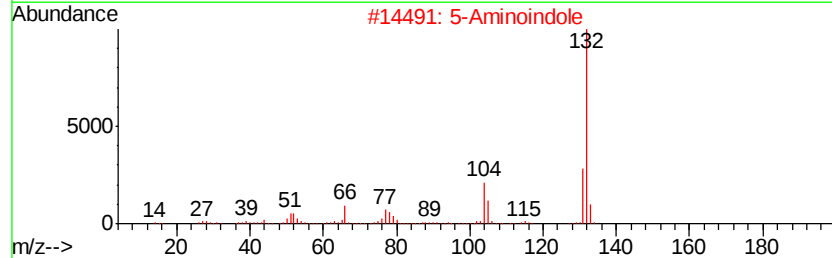
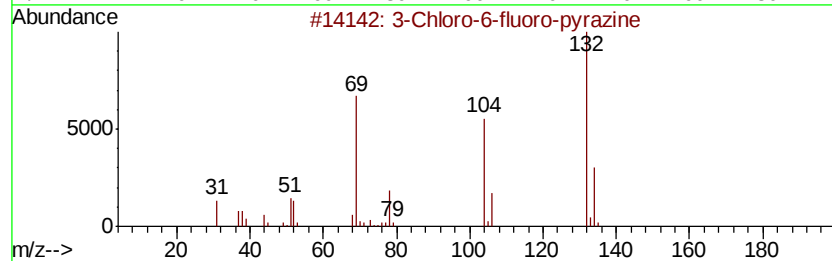
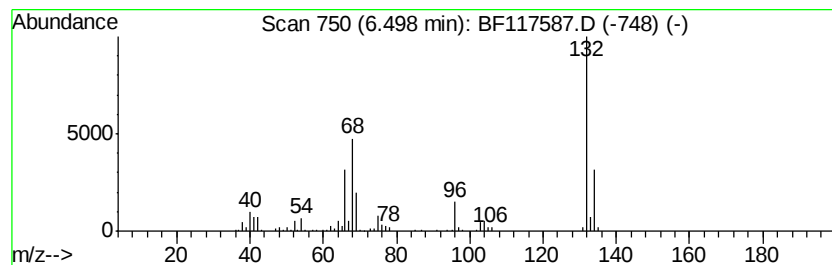
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	56.57 ng	749651	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Xenon	132	Xe	007440-63-3	9



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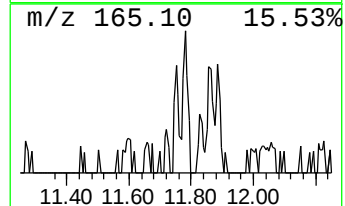
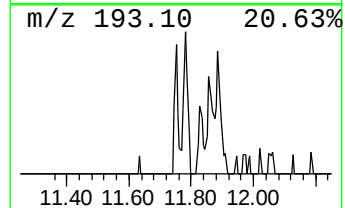
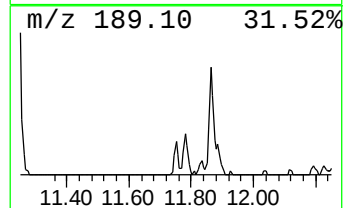
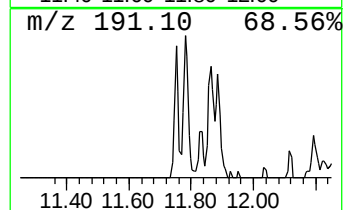
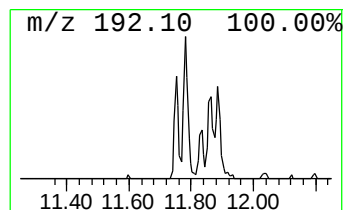
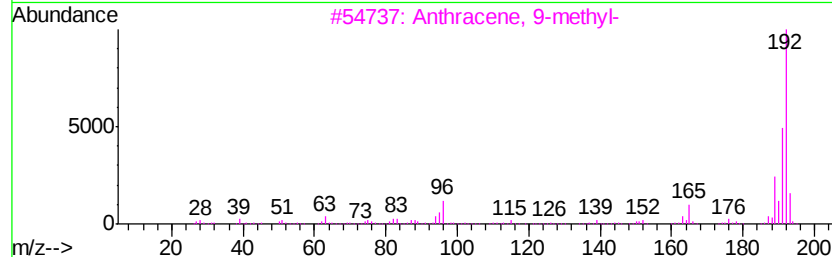
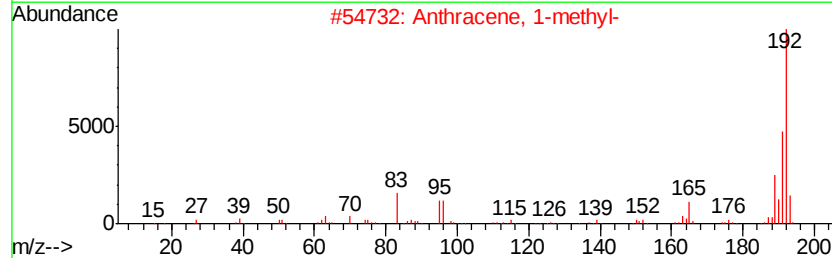
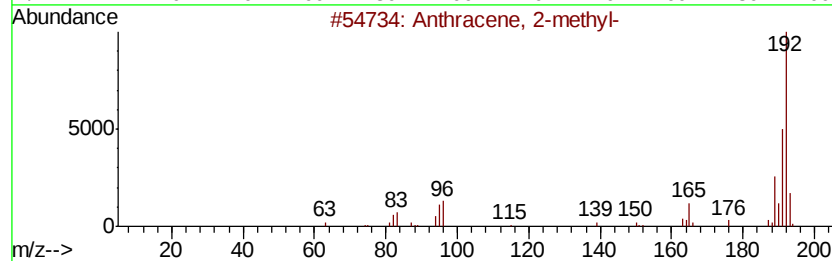
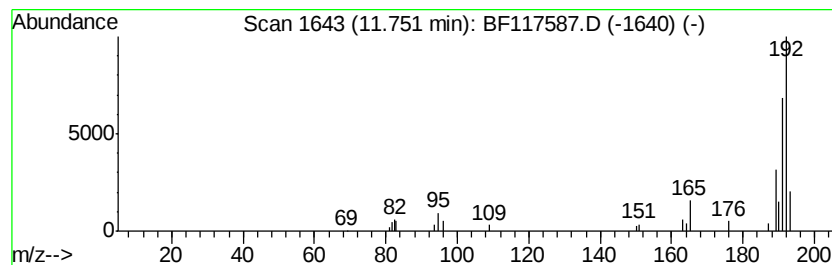
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	2.08 ng	28345	Phenanthrene-d10	11.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
4			1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	90
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87



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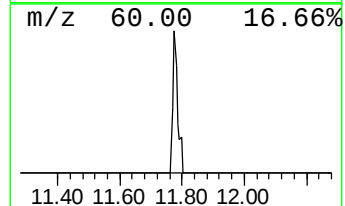
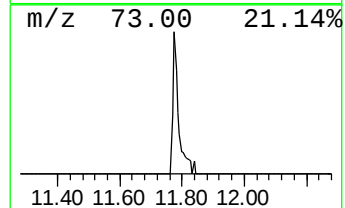
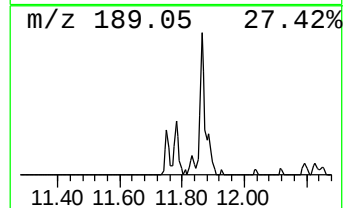
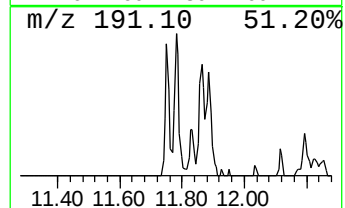
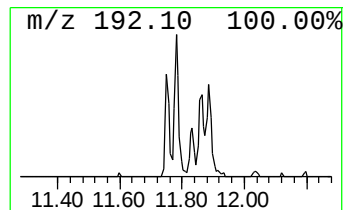
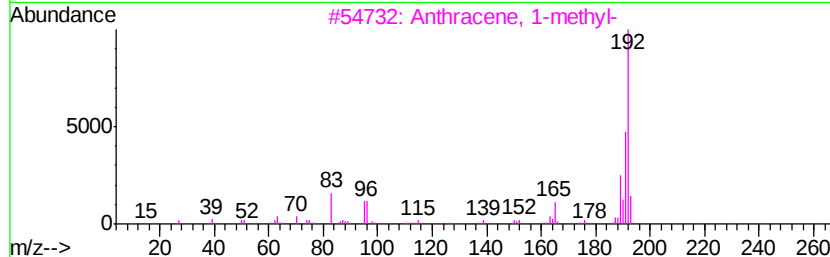
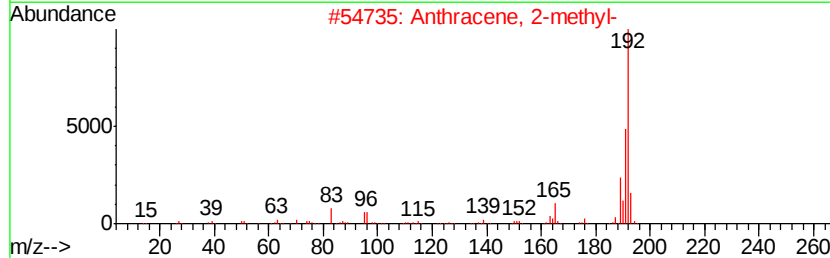
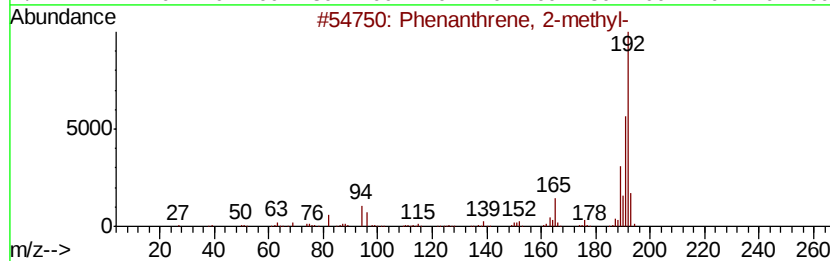
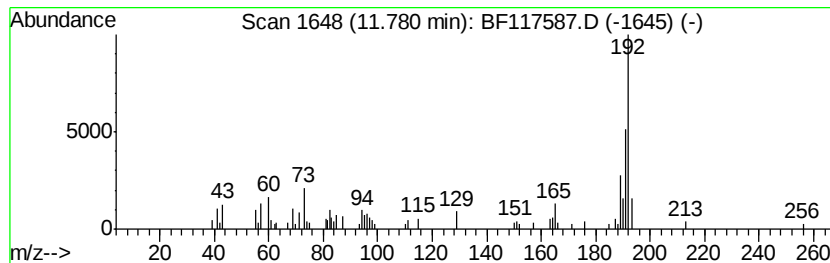
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	6.12 ng	83324	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
5		1H-Cyclopropa[1]phenanthrene, 1a,...	192	C15H12	000949-41-7	92



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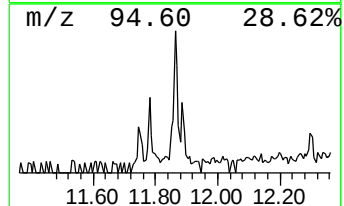
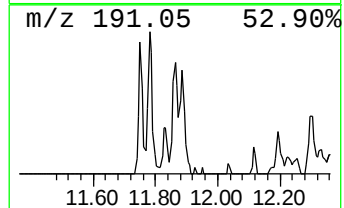
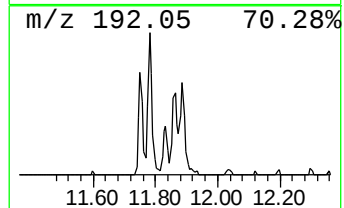
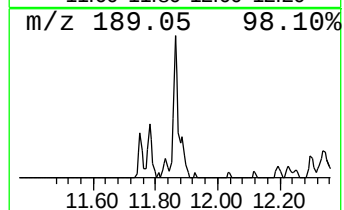
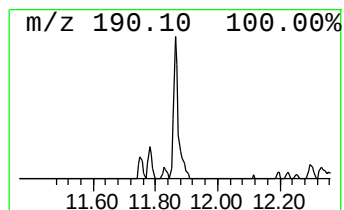
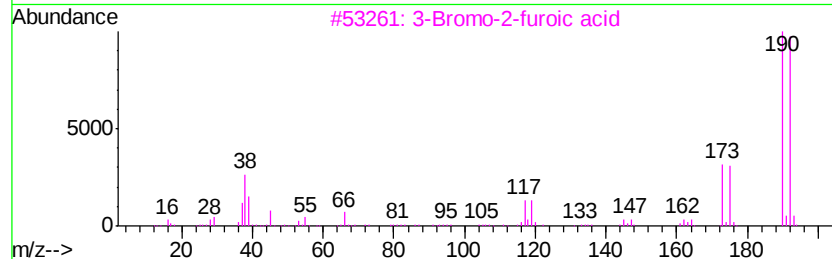
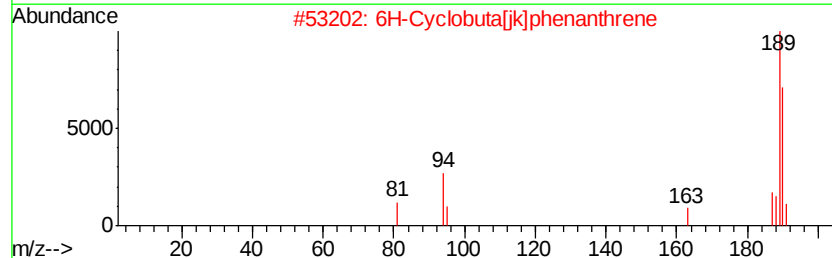
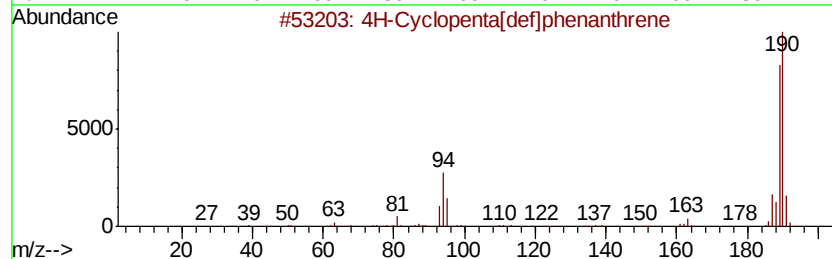
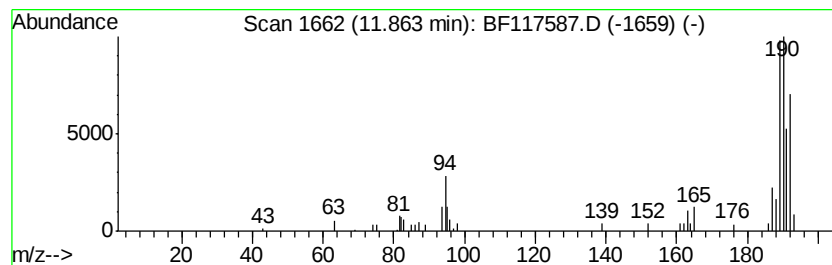
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.86	4.22 ng	57520	Phenanthrene-d10		11.24	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	62
3		3-Bromo-2-furoic acid	190	C5H3BrO3	014903-90-3	27
4		2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	25
5		Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
Data File : BF117587.D
Acq On : 29 Oct 2019 13:37
Operator : JU
Sample : K5503-13
Misc :
ALS Vial : 10 Sample Multiplier: 1

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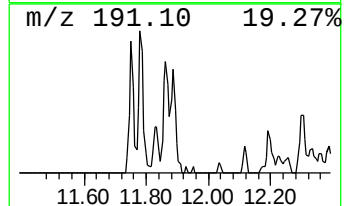
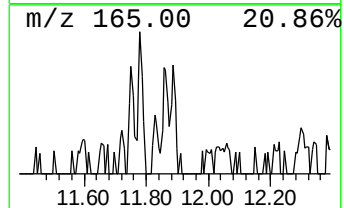
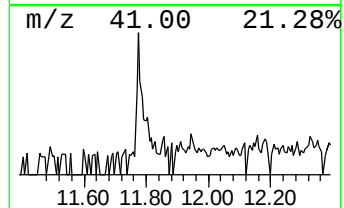
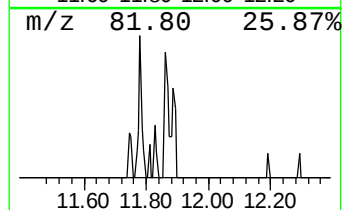
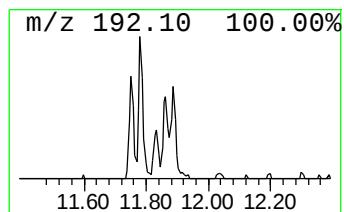
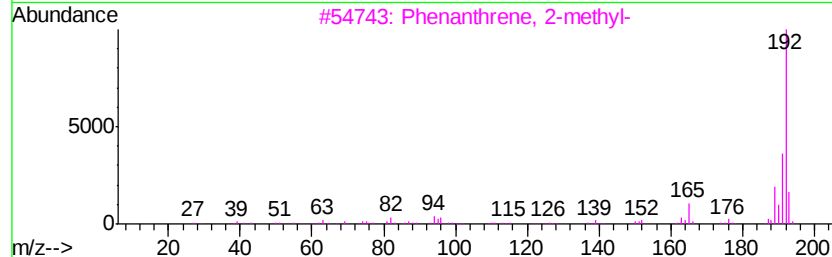
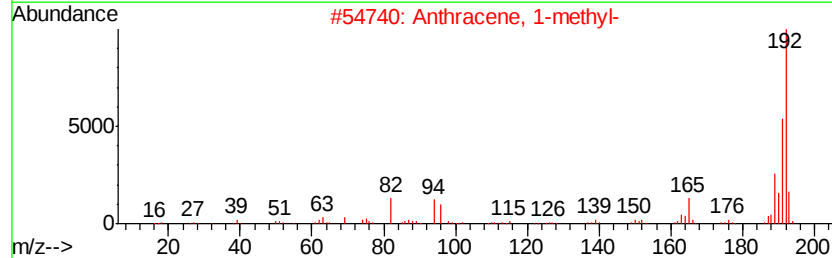
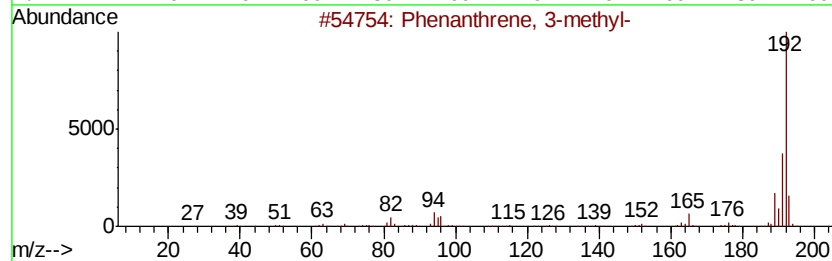
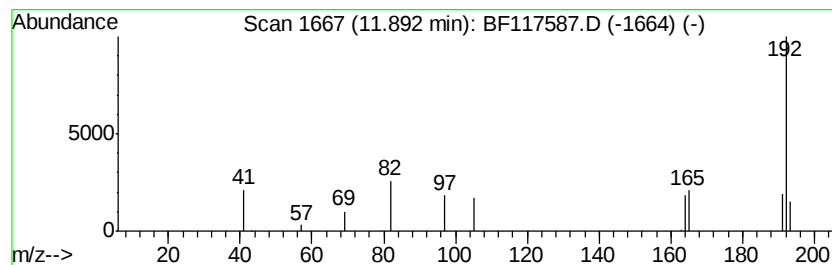
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	2.10 ng	28605	Phenanthrene-d10	11.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	90
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	86
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	78
5		1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
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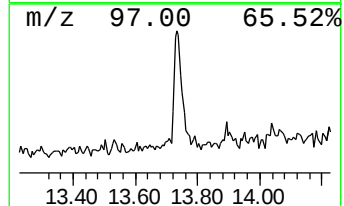
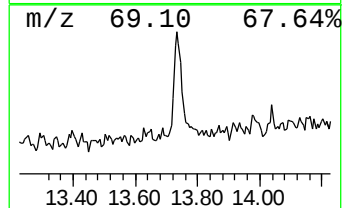
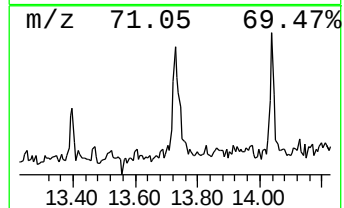
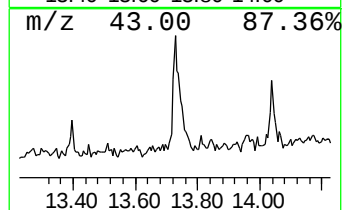
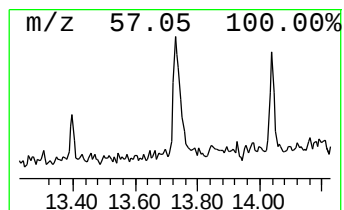
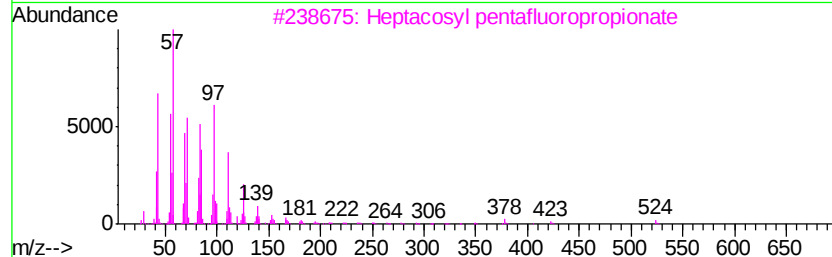
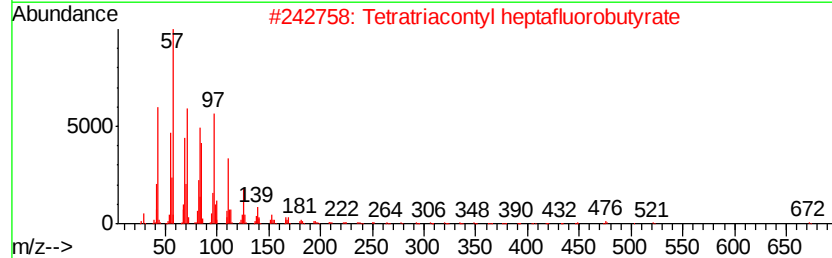
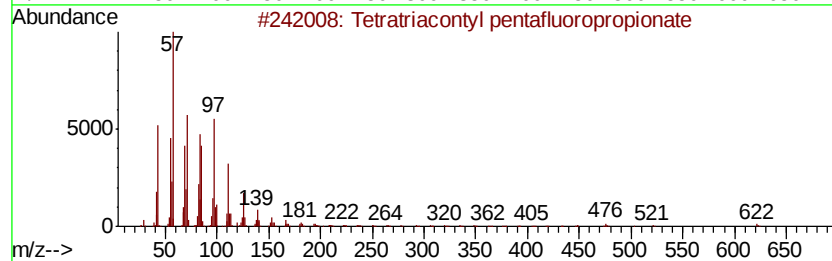
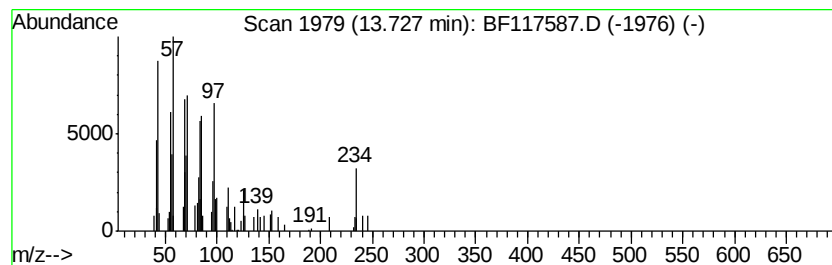
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tetratriacontyl pentafluoro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	2.40 ng	56002	Chrysene-d12	13.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	81
2		Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	81
3		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	68
4		Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	68
5		17-Pentatriacontene	491	C35H70	006971-40-0	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
Data File : BF117587.D
Acq On : 29 Oct 2019 13:37
Operator : JU
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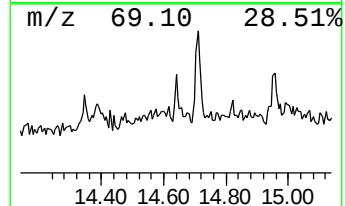
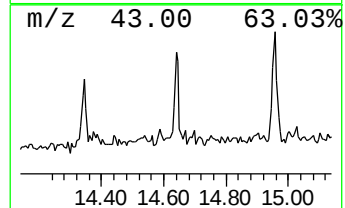
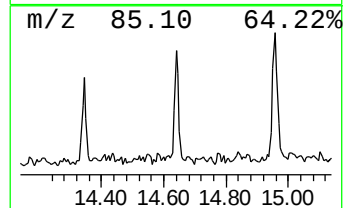
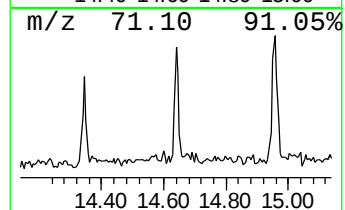
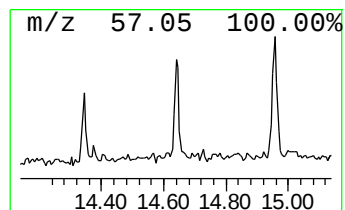
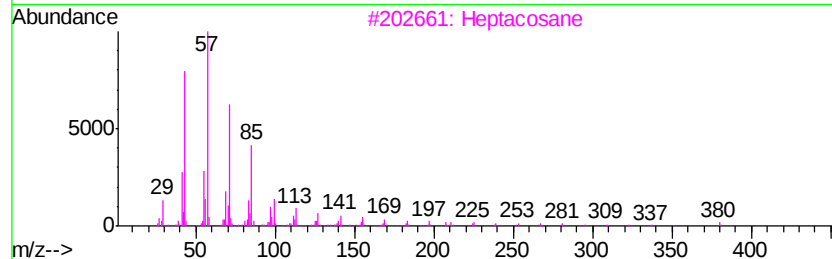
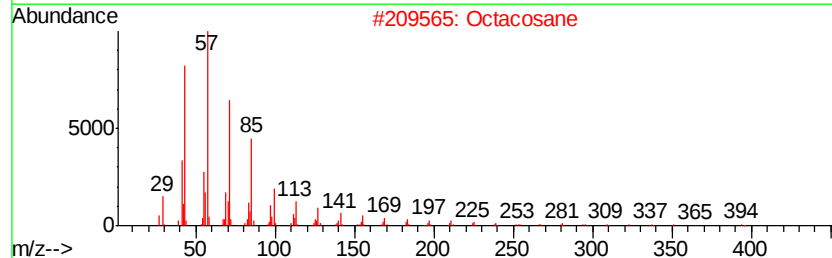
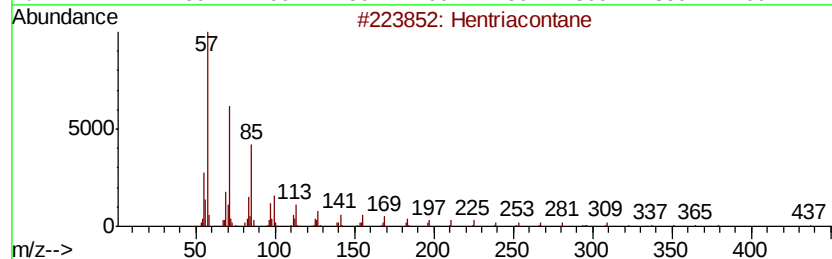
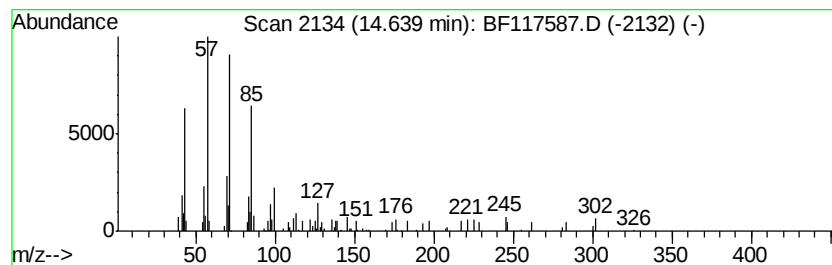
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hentriacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	3.39 ng	39419	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	80
2		Octacosane	394	C28H58	000630-02-4	80
3		Heptacosane	380	C27H56	000593-49-7	72
4		6,6-Diethylhoctadecane	310	C22H46	1000360-41-8	72
5		Tetratetracontane	619	C44H90	007098-22-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
Data File : BF117587.D
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Misc :
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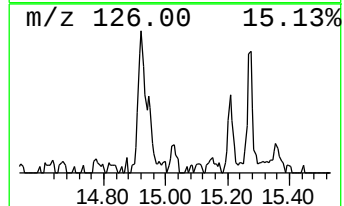
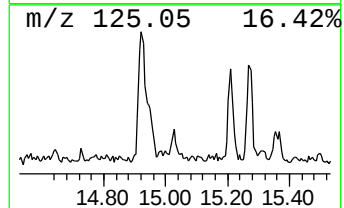
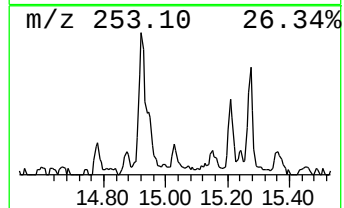
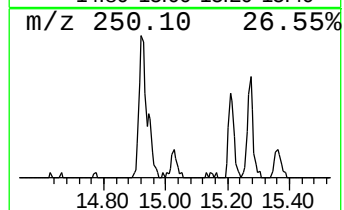
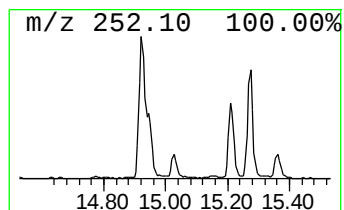
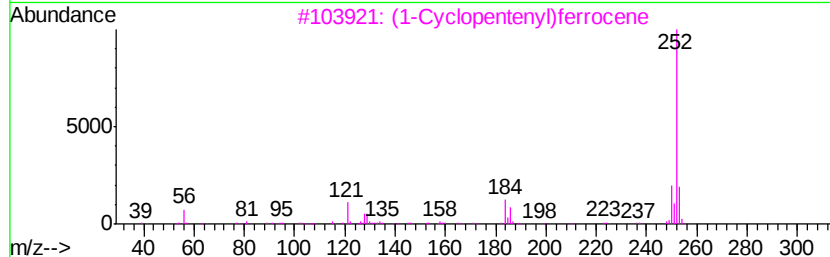
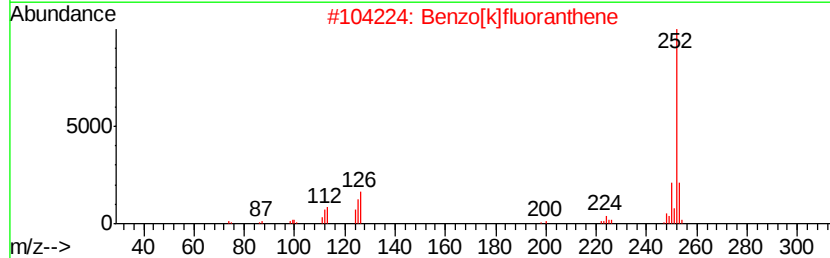
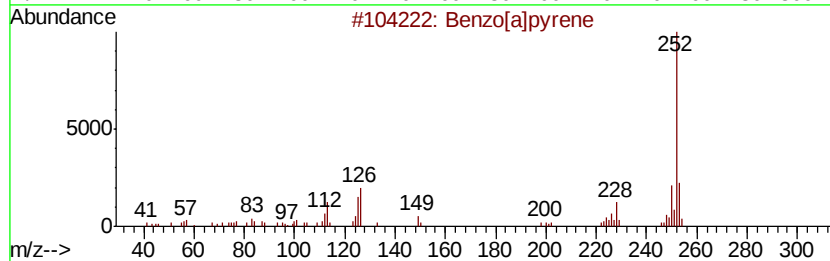
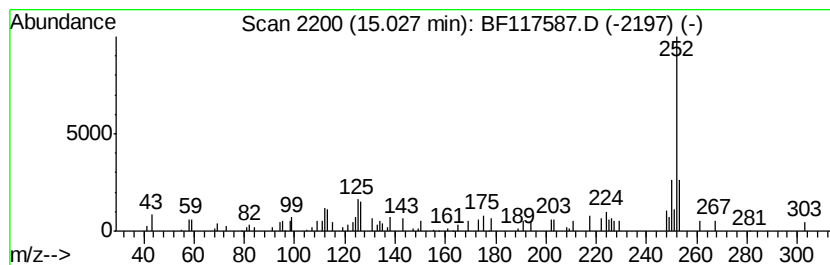
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.03	2.55 ng	29604	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[a]pyrene	252	C20H12	000050-32-8	70
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	70
3		(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	64
4		Perylene	252	C20H12	000198-55-0	62
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
Data File : BF117587.D
Acq On : 29 Oct 2019 13:37
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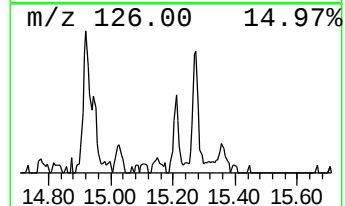
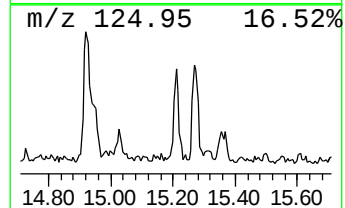
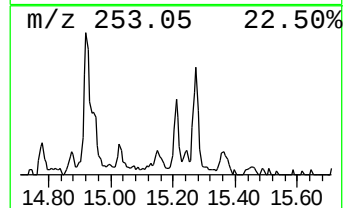
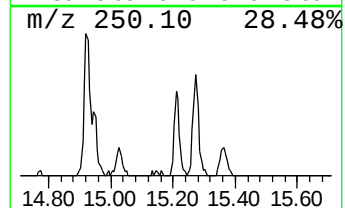
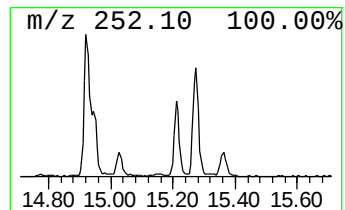
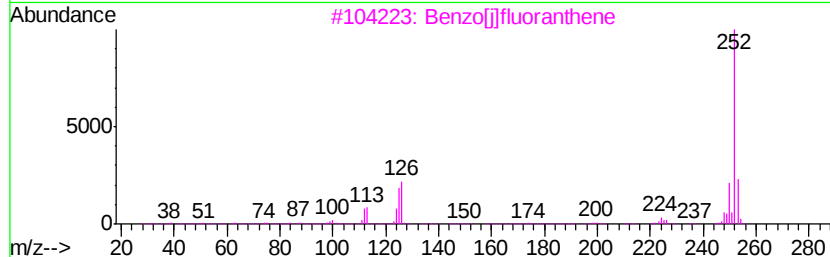
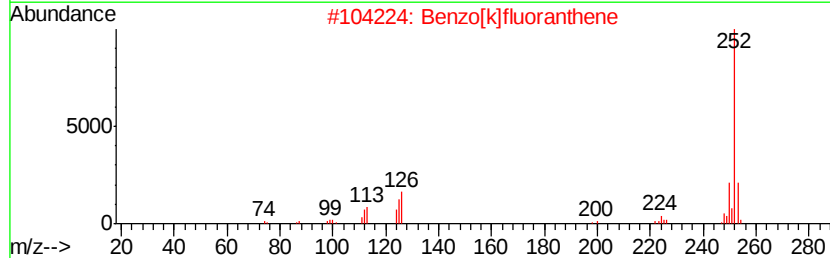
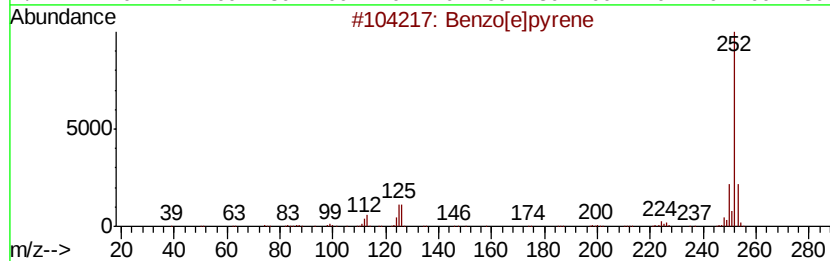
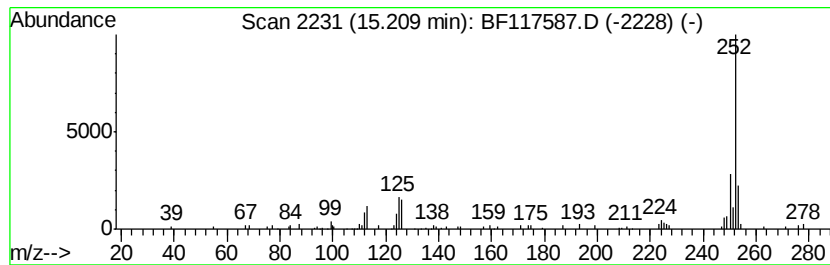
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	5.48 ng	63709	Perylene-d12	15.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	96
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3			Benzo[i]fluoranthene	252	C20H12	000205-82-3	95
4			Perylene	252	C20H12	000198-55-0	94
5			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
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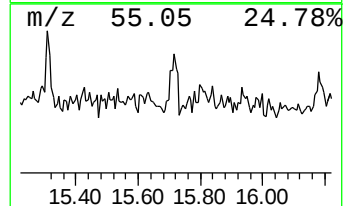
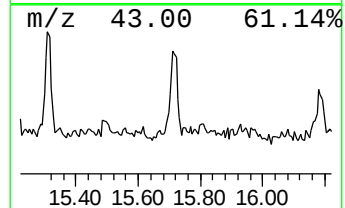
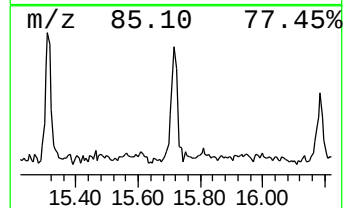
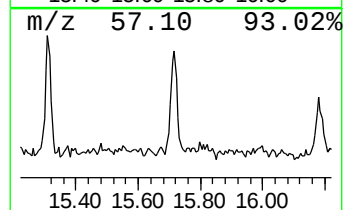
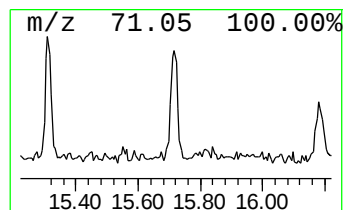
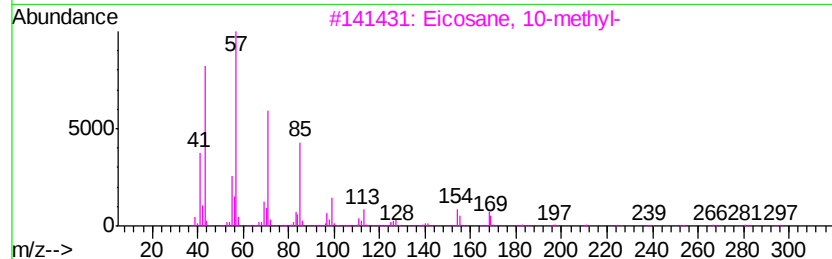
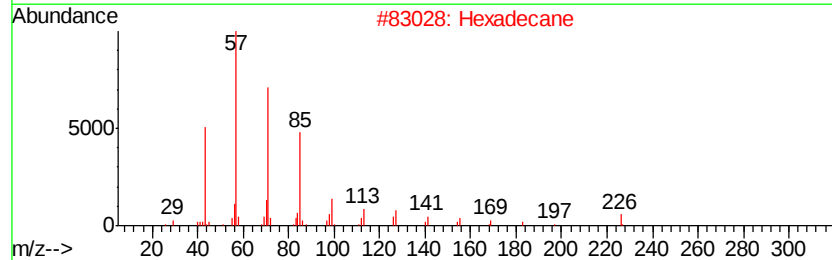
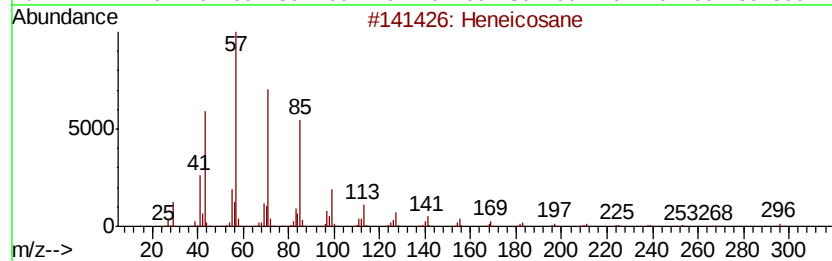
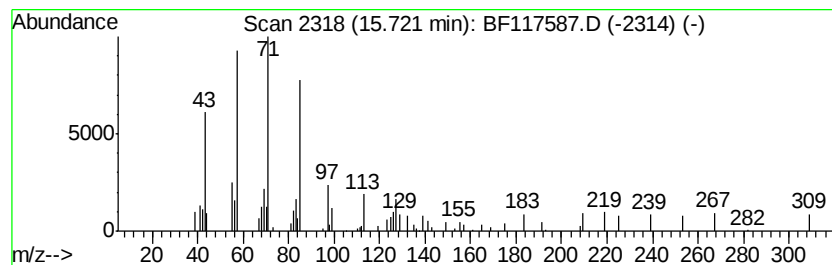
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	3.68 ng	42726	Perylene-d12	15.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	64
2			Hexadecane	226	C16H34	000544-76-3	59
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	59
4			Octacosane	394	C28H58	000630-02-4	58
5			Hentriacontane	437	C31H64	000630-04-6	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
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ClientSampleId :
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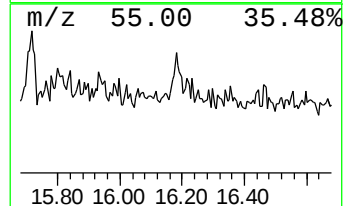
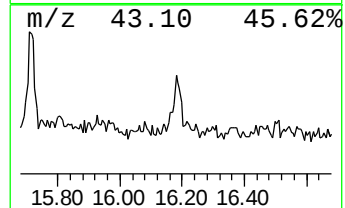
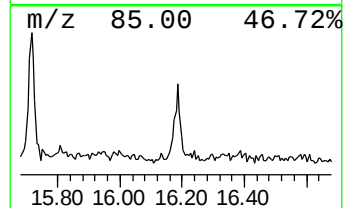
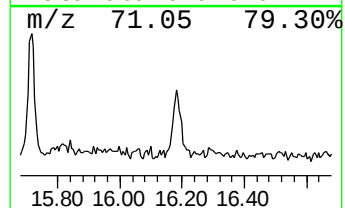
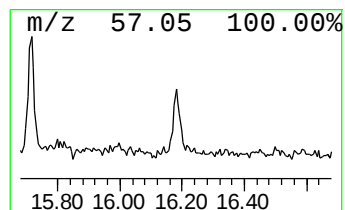
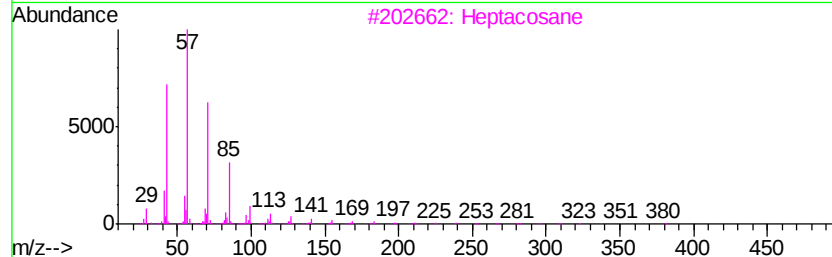
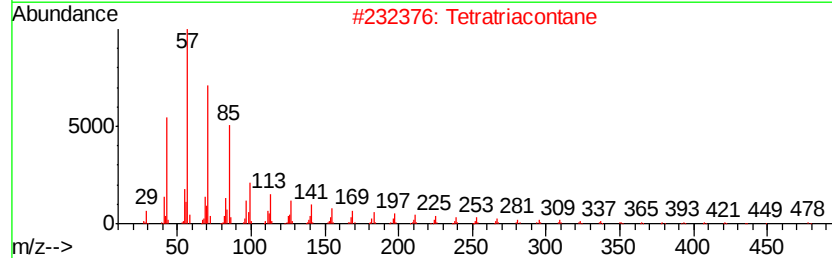
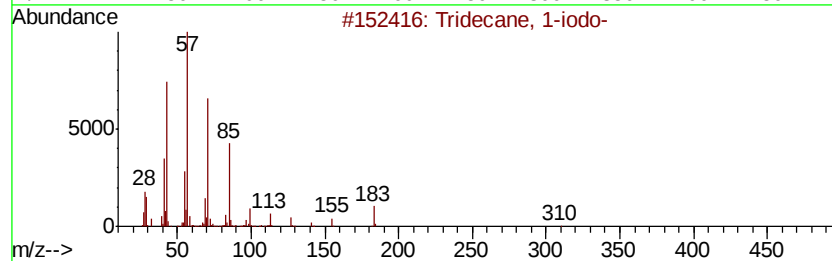
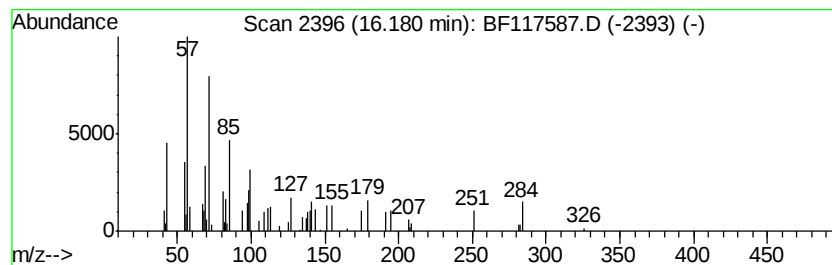
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.18 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	2.41 ng	28020	Perylene-d12	15.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	47
2			Tetratriacontane	479	C34H70	014167-59-0	47
3			Heptacosane	380	C27H56	000593-49-7	47
4			Eicosane	282	C20H42	000112-95-8	47
5			5,5,7,7-Tetraethylundecane	268	C19H40	1000360-42-0	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102919\
 Data File : BF117587.D
 Acq On : 29 Oct 2019 13:37
 Operator : JU
 Sample : K5503-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 172-80-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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
unknown6.50	6.50	56.6	ng	749651	1	6.73	265020	20.0
Anthracene, 2-met...	11.75	2.1	ng	28345	4	11.24	272472	20.0
Phenanthrene, 2-m...	11.78	6.1	ng	83324	4	11.24	272472	20.0
4H-Cyclopenta[def...	11.86	4.2	ng	57520	4	11.24	272472	20.0
Phenanthrene, 3-m...	11.89	2.1	ng	28605	4	11.24	272472	20.0
Tetratriacontyl p...	13.73	2.4	ng	56002	5	13.89	466360	20.0
Hentriacontane	14.64	3.4	ng	39419	6	15.33	232312	20.0
Perylene	15.03	2.5	ng	29604	6	15.33	232312	20.0
Benzo[e]pyrene	15.21	5.5	ng	63709	6	15.33	232312	20.0
Heneicosane	15.72	3.7	ng	42726	6	15.33	232312	20.0
unknown16.18	16.18	2.4	ng	28020	6	15.33	232312	20.0