

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082019\
 Data File : BF116455.D
 Acq On : 21 Aug 2019 2:42
 Operator : HP/JU
 Sample : K4354-02 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PX-02-081319

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-BF073019.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	3.810	289	293	309	rBV	124236	222740	10.41%	0.781%
2	5.440	567	570	609	rBV	425153	879305	41.10%	3.084%
3	6.451	740	742	761	rBV	603604	970162	45.35%	3.402%
4	6.581	761	764	778	rVB	964791	1053869	49.26%	3.696%
5	6.804	799	802	813	rBV	566973	549384	25.68%	1.927%
6	6.957	825	828	845	rBV	818088	797802	37.29%	2.798%
7	7.369	895	898	927	rBV	457974	642897	30.05%	2.255%
8	8.081	1017	1019	1044	rVB	666375	775838	36.27%	2.721%
9	9.151	1198	1201	1213	rBV	1307064	1181104	55.21%	4.142%
10	9.551	1266	1269	1277	rVB	176984	160594	7.51%	0.563%
11	9.833	1314	1317	1321	rBV	1086973	920748	43.04%	3.229%
12	9.869	1321	1323	1331	rVB	147577	138077	6.45%	0.484%
13	10.051	1349	1354	1363	rBV	37234	46517	2.17%	0.163%
14	10.386	1408	1411	1418	rBV	79221	85320	3.99%	0.299%
15	10.628	1449	1452	1466	rBV	766851	838523	39.20%	2.941%
16	11.151	1538	1541	1543	rBV	35269	36258	1.69%	0.127%
17	11.175	1543	1545	1549	rVV2	30127	29641	1.39%	0.104%
18	11.222	1551	1553	1559	rVB	65018	60463	2.83%	0.212%
19	11.322	1566	1570	1572	rBV	1260719	1086735	50.80%	3.811%
20	11.345	1572	1574	1579	rVV	1606945	1339119	62.59%	4.696%
21	11.404	1579	1584	1591	rVB	281190	311978	14.58%	1.094%
22	11.563	1608	1611	1617	rBV	178939	190060	8.88%	0.667%
23	11.827	1653	1656	1658	rBV	113926	108853	5.09%	0.382%
24	11.857	1658	1661	1667	rVV	154361	171752	8.03%	0.602%
25	11.904	1667	1669	1672	rVV	48140	45348	2.12%	0.159%
26	11.939	1672	1675	1677	rVV	224514	225987	10.56%	0.793%
27	11.963	1677	1679	1684	rVB	68140	64964	3.04%	0.228%
28	12.116	1703	1705	1710	rVB	104540	93388	4.37%	0.328%
29	12.169	1710	1714	1721	rBV2	82059	105622	4.94%	0.370%
30	12.374	1745	1749	1751	rBV	61728	59441	2.78%	0.208%
31	12.474	1761	1766	1773	rVB2	71400	104324	4.88%	0.366%
32	12.539	1773	1777	1782	rBV	2309528	2139351	100.00%	7.503%
33	12.616	1785	1790	1795	rVV2	25931	45324	2.12%	0.159%
34	12.692	1799	1803	1806	rBV	108447	105406	4.93%	0.370%

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 Sampling : 1
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35	12.763	1809	1815	1821	rVV	1807775	2122756	99.22%	7.445%
36	12.822	1821	1825	1828	rVB	73321	84267	3.94%	0.296%
37	12.898	1834	1838	1843	rBV	1709895	1545473	72.24%	5.420%
38	12.939	1843	1845	1849	rVB	88826	91459	4.28%	0.321%
39	12.992	1849	1854	1857	rBV2	83527	142680	6.67%	0.500%
40	13.027	1857	1860	1864	rVB	44396	42648	1.99%	0.150%
41	13.074	1864	1868	1869	rBV3	70357	80611	3.77%	0.283%
42	13.098	1869	1872	1877	rVB	192971	210348	9.83%	0.738%
43	13.163	1880	1883	1885	rVV	151244	131187	6.13%	0.460%
44	13.198	1885	1889	1896	rVV3	124217	228893	10.70%	0.803%
45	13.257	1897	1899	1902	rVB3	31919	30446	1.42%	0.107%
46	13.292	1902	1905	1908	rBV	64090	67118	3.14%	0.235%
47	13.327	1908	1911	1920	rVB3	71837	109145	5.10%	0.383%
48	13.463	1930	1934	1937	rBV5	17704	30110	1.41%	0.106%
49	13.621	1957	1961	1966	rVB	96328	111294	5.20%	0.390%
50	13.721	1974	1978	1979	rBV	135176	146121	6.83%	0.512%
51	13.769	1984	1986	1988	rVV	124028	130790	6.11%	0.459%
52	13.827	1992	1996	2001	rVV2	145284	205216	9.59%	0.720%
53	13.886	2004	2006	2009	rVV	53237	53407	2.50%	0.187%
54	13.921	2009	2012	2013	rVV	110111	99078	4.63%	0.347%
55	13.963	2014	2019	2022	rVV2	1281077	1978009	92.46%	6.937%
56	13.992	2022	2024	2031	rVB	778617	721686	33.73%	2.531%
57	14.074	2035	2038	2041	rVV	139366	128078	5.99%	0.449%
58	14.133	2046	2048	2052	rBV2	29459	30126	1.41%	0.106%
59	14.168	2052	2054	2056	rBV	38843	32686	1.53%	0.115%
60	14.357	2081	2086	2089	rBV	109291	132727	6.20%	0.465%
61	14.404	2089	2094	2096	rBV3	53196	71832	3.36%	0.252%
62	14.427	2096	2098	2100	rVB2	41305	31307	1.46%	0.110%
63	14.457	2100	2103	2105	rBV2	44314	42636	1.99%	0.150%
64	14.480	2105	2107	2109	rBV2	52473	56164	2.63%	0.197%
65	14.704	2142	2145	2148	rVB3	87984	101199	4.73%	0.355%
66	15.004	2192	2196	2199	rBV	678556	990412	46.29%	3.473%
67	15.115	2212	2215	2222	rVB	114990	135643	6.34%	0.476%
68	15.298	2243	2246	2251	rBV	353768	414304	19.37%	1.453%
69	15.363	2254	2257	2264	rBV	548110	662207	30.95%	2.322%
70	15.421	2264	2267	2271	rBV	612696	713746	33.36%	2.503%
71	16.892	2512	2517	2527	rVB	223721	480402	22.46%	1.685%

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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72	17.062	2543	2546	2551	rVB	30894	47966	2.24%	0.168%
73	17.333	2587	2592	2602	rVB	171469	389435	18.20%	1.366%
74	17.609	2635	2639	2649	rVB	44352	103064	4.82%	0.361%

Sum of corrected areas: 28513570

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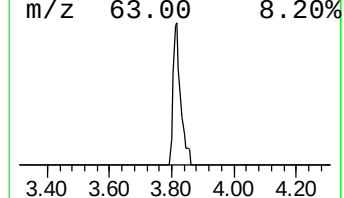
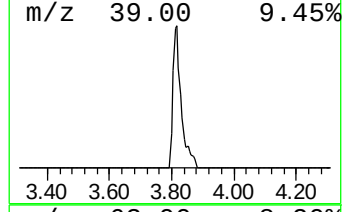
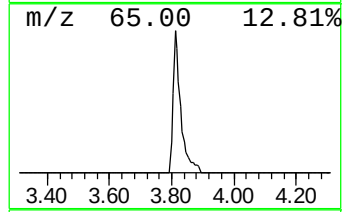
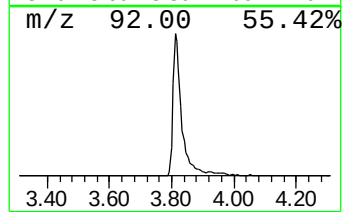
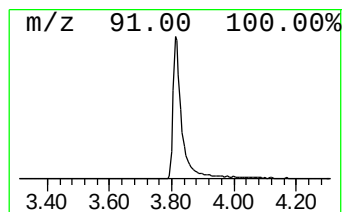
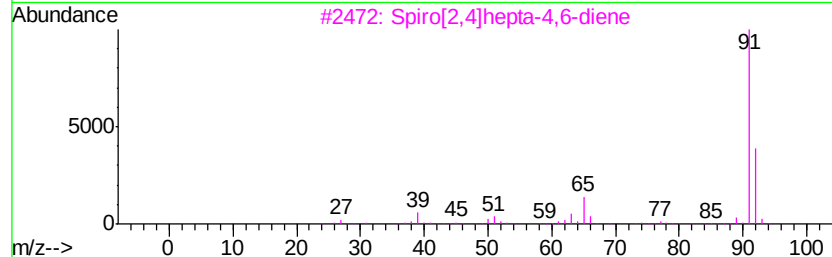
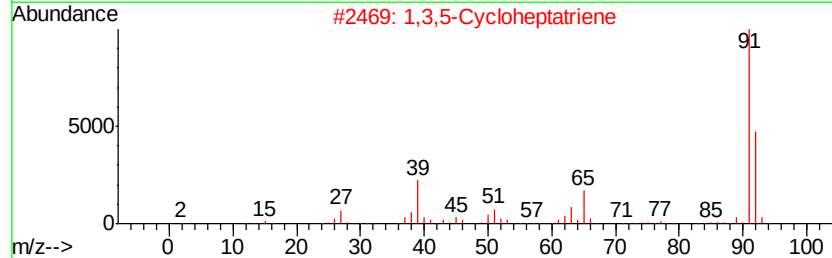
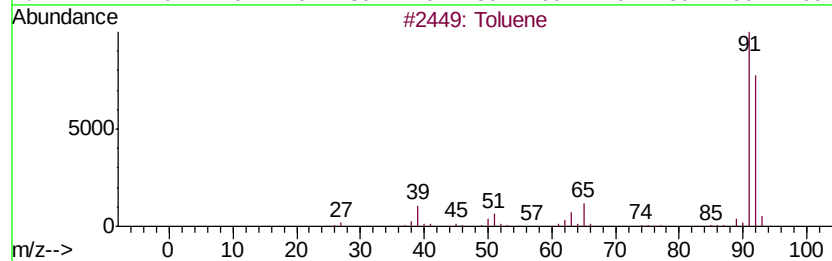
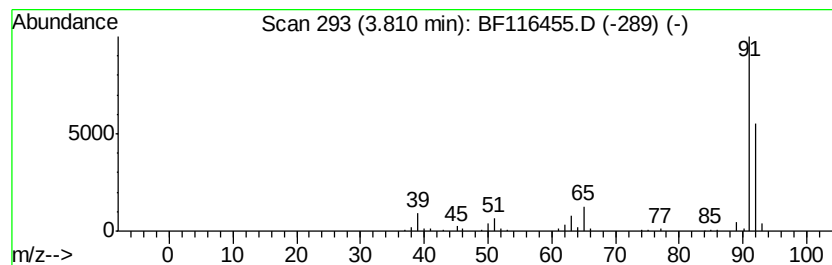
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1,3,5-Cycloheptatriene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.81	8.11 ng	222740	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	95
2			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90
3			Spiro[2.4]hepta-4,6-diene	92	C7H8	000765-46-8	87
4			2,5-Norbornadiene	92	C7H8	000121-46-0	64
5			Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	59



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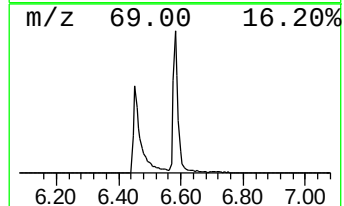
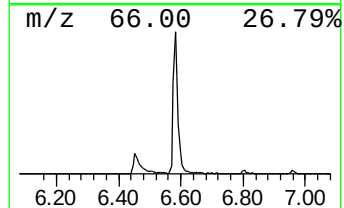
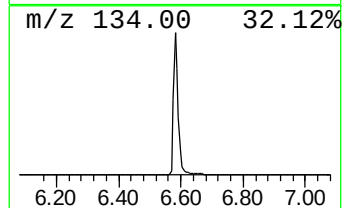
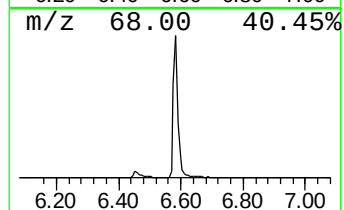
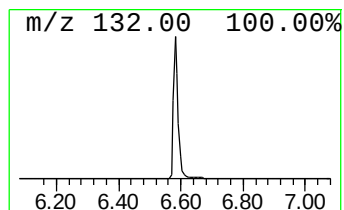
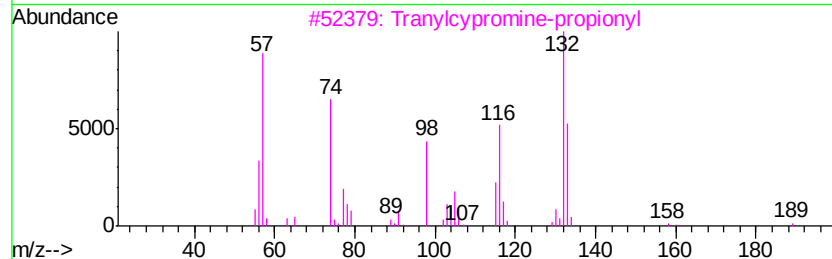
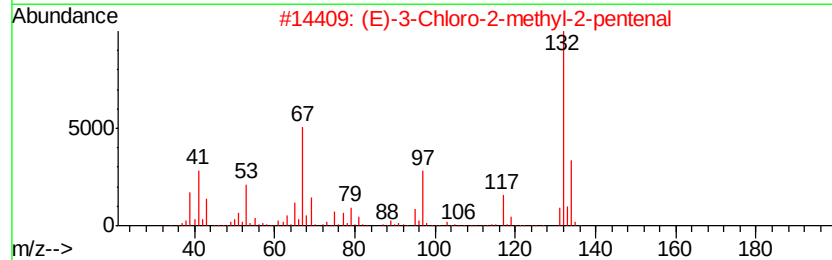
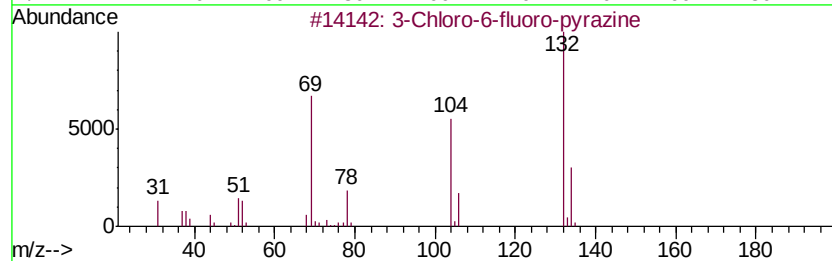
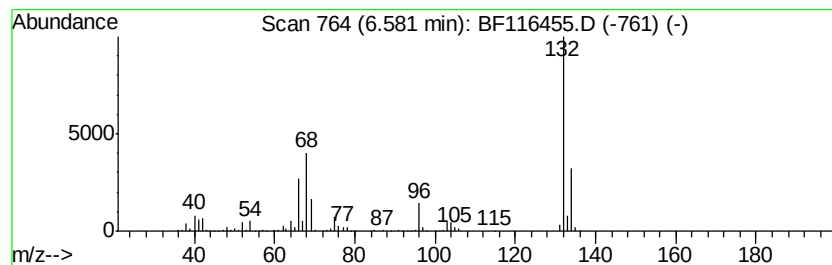
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	38.37 ng	1053870	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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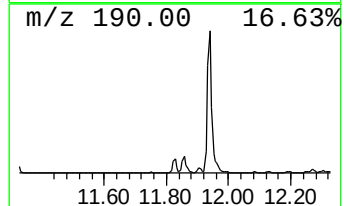
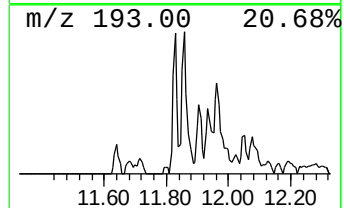
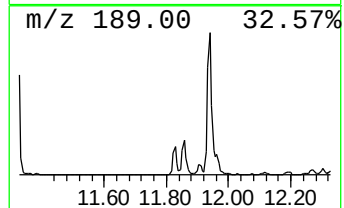
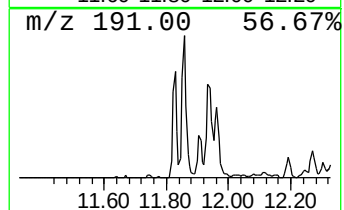
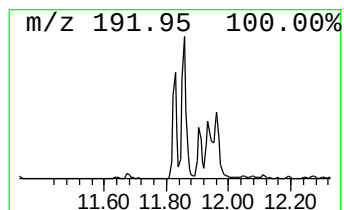
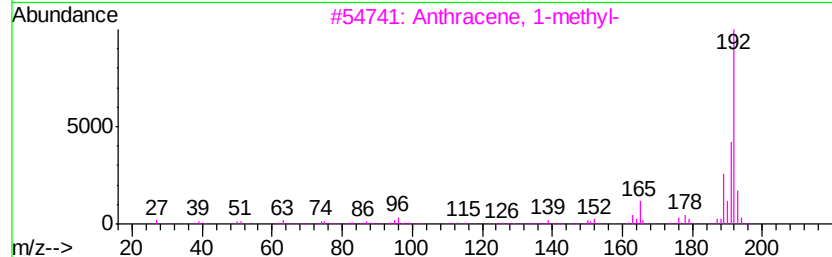
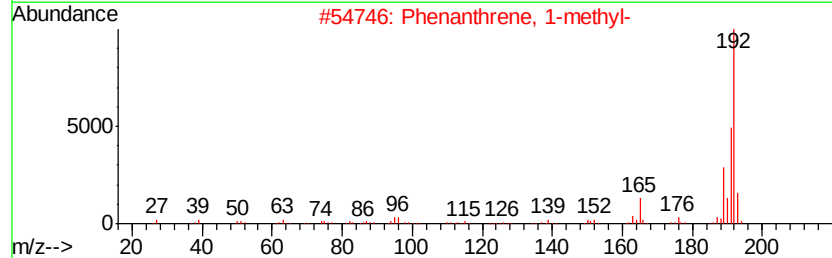
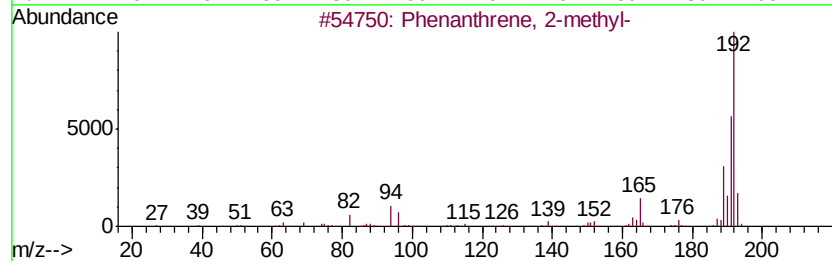
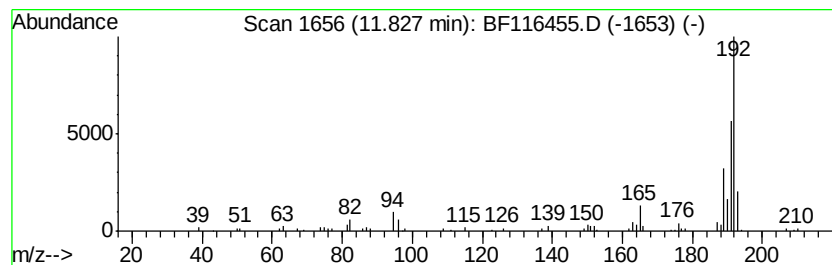
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	2.00 ng	108853	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91



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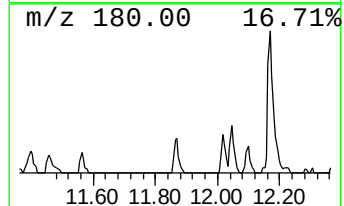
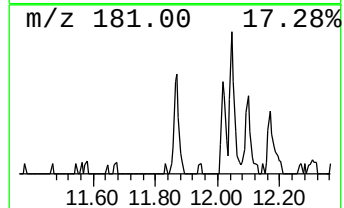
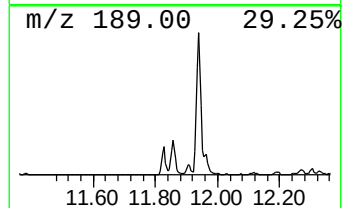
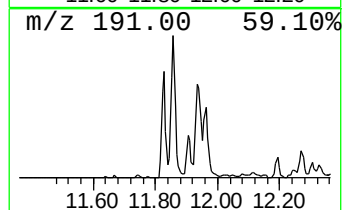
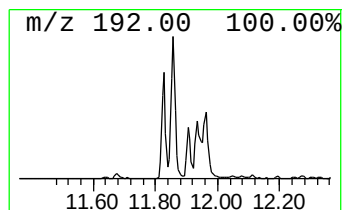
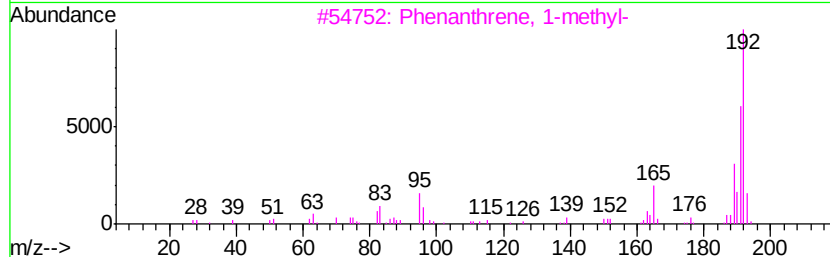
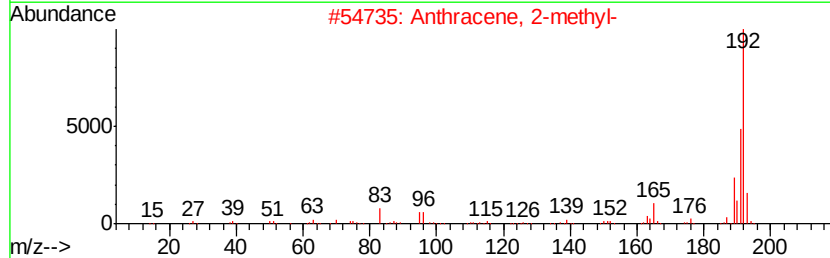
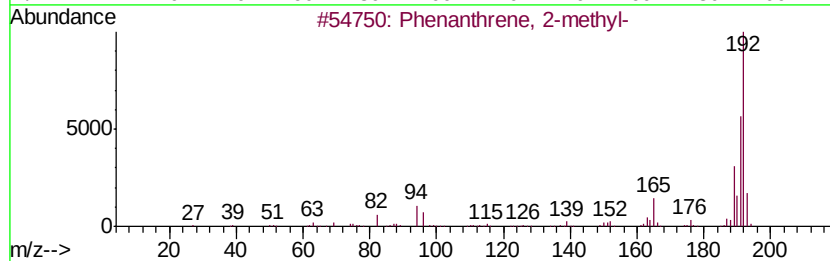
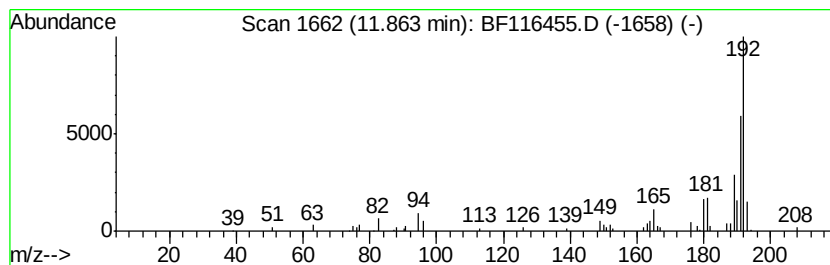
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.86	3.16 ng	171752	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
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Operator : HP/JU
Sample : K4354-02 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

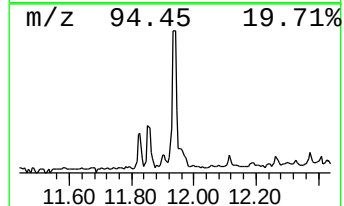
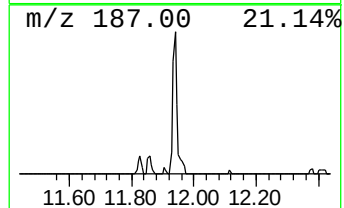
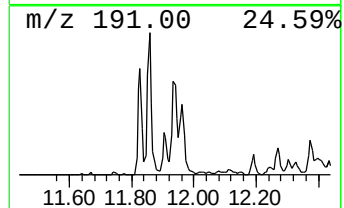
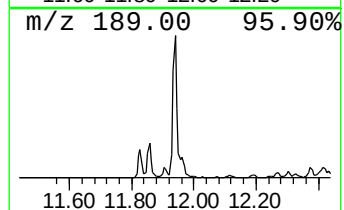
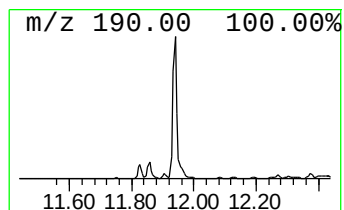
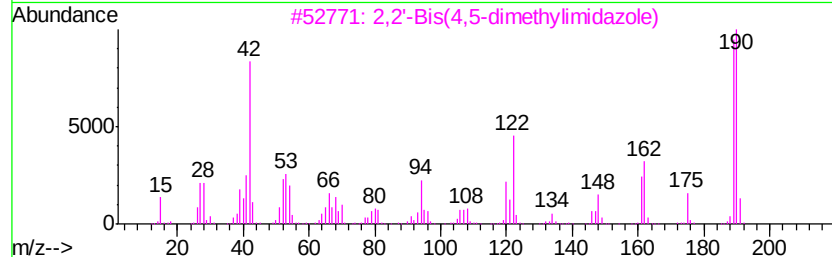
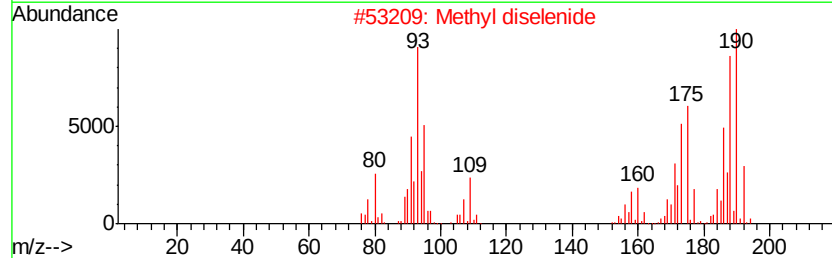
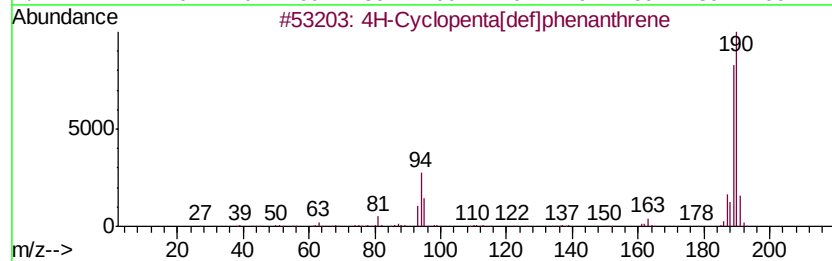
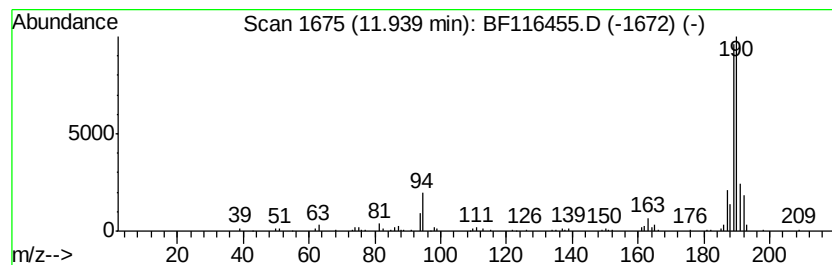
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.94	4.16 ng	225987	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	Methyl diselenide	190	C2H6Se2	007101-31-7	55
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
4	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
Data File : BF116455.D
Acq On : 21 Aug 2019 2:42
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Sample : K4354-02 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

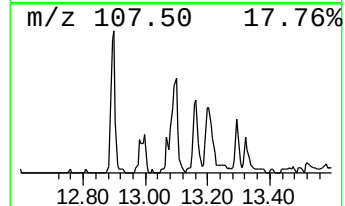
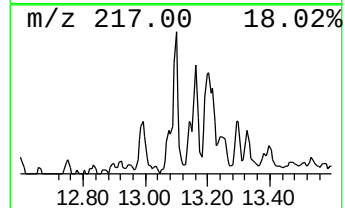
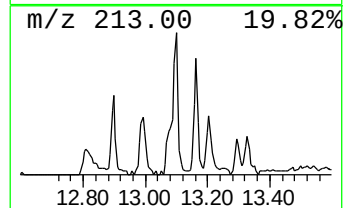
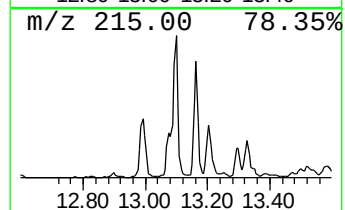
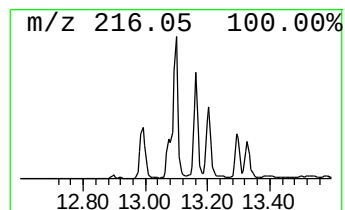
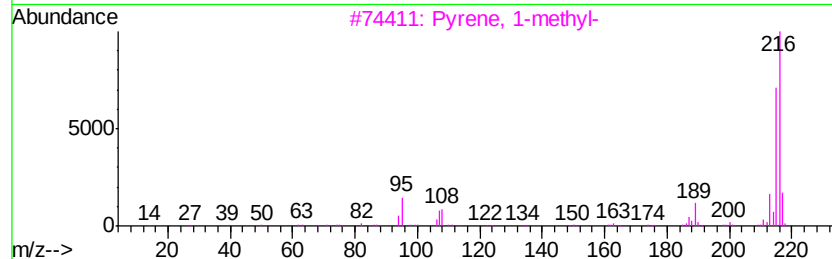
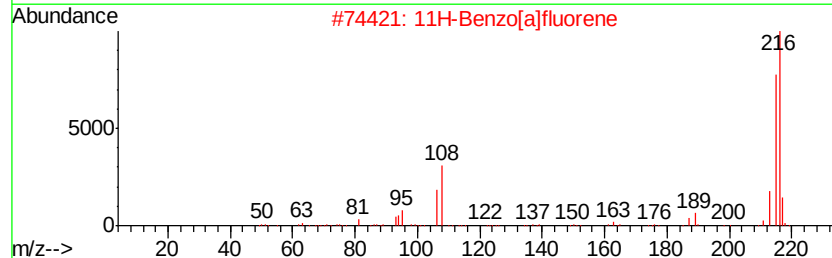
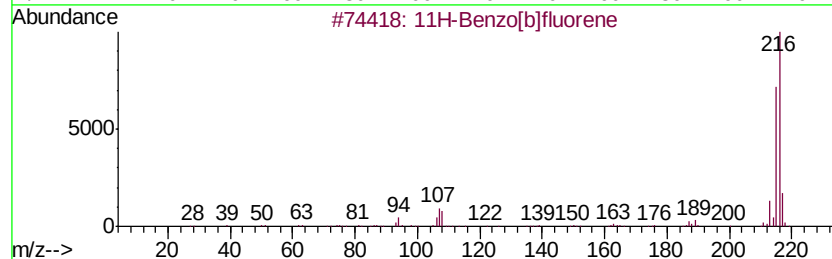
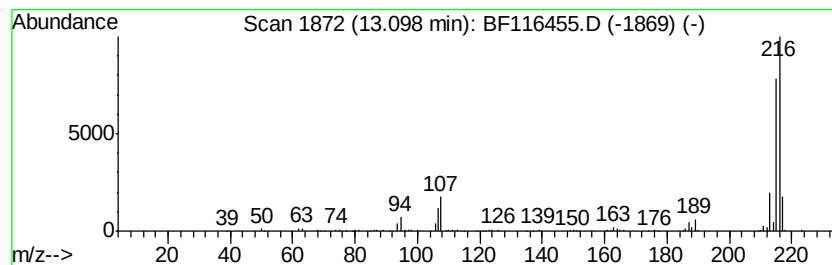
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.10	2.13 ng	210348	Chrysene-d12	13.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	86
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
Data File : BF116455.D
Acq On : 21 Aug 2019 2:42
Operator : HP/JU
Sample : K4354-02 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

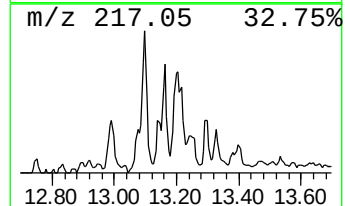
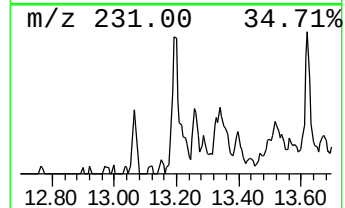
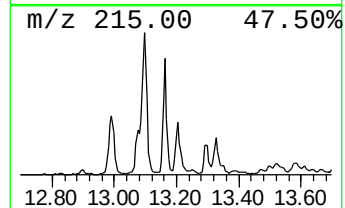
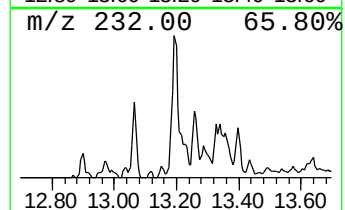
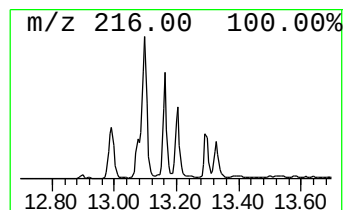
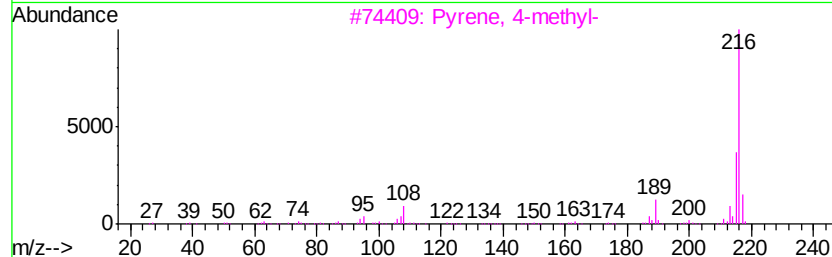
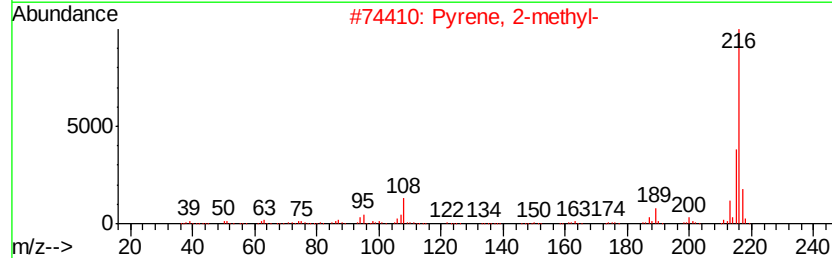
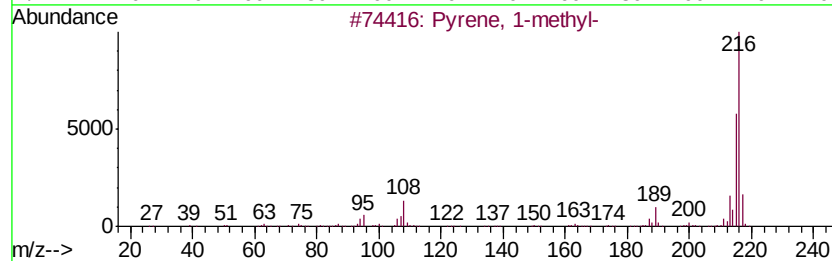
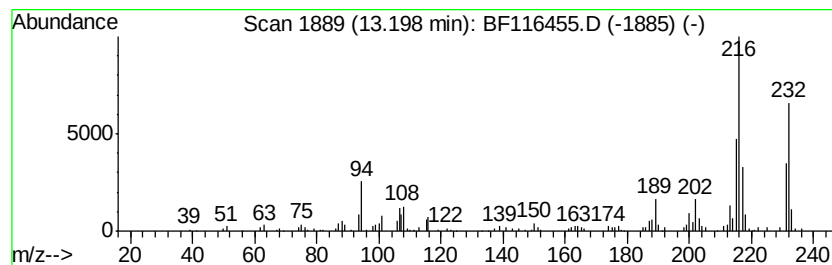
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.20	2.31 ng	228893	Chrysene-d12	13.96	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	55
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	55
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	55
4	1-benzyl-3-methylnaphthalene	232	C18H16	1000379-93-5	46
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
Data File : BF116455.D
Acq On : 21 Aug 2019 2:42
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Sample : K4354-02 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

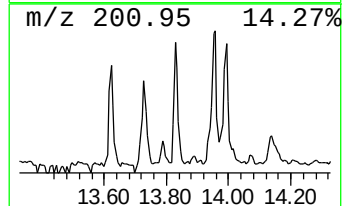
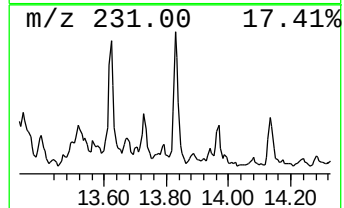
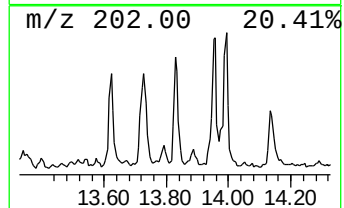
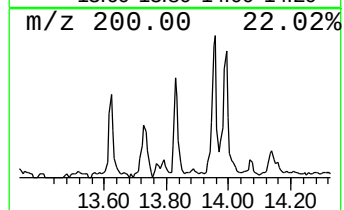
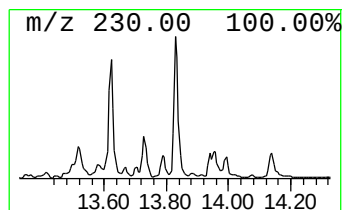
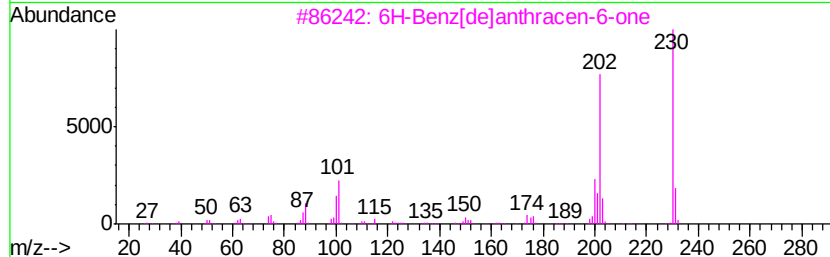
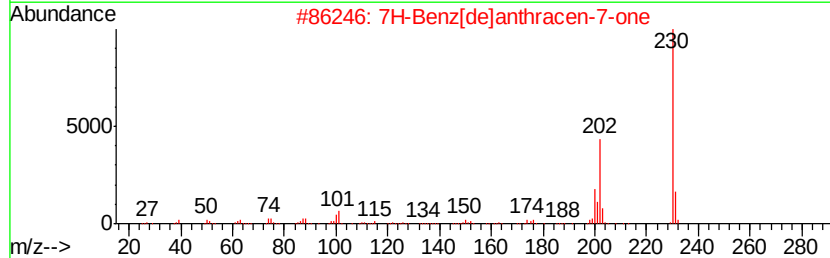
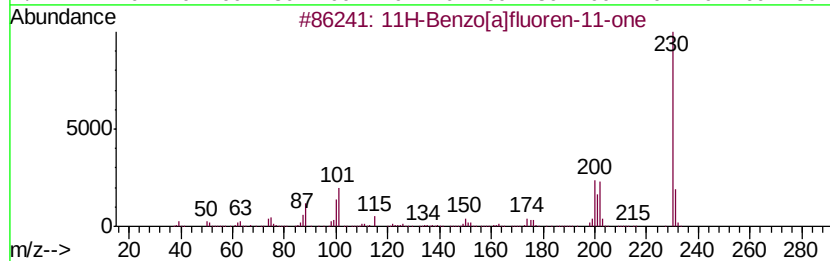
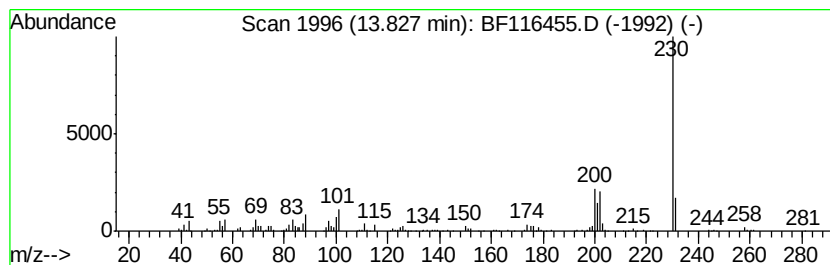
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7H-Benz[de]anthracen-7-one Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.07 ng	205216	Chrysene-d12	13.96
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[a]fluoren-11-one	230 C17H10O	000479-79-8 96
2		7H-Benz[de]anthracen-7-one	230 C17H10O	000082-05-3 90
3		6H-Benz[de]anthracen-6-one	230 C17H10O	080252-14-8 64
4		1-Pvrenecarboxaldehyde	230 C17H10O	003029-19-4 64
5		o-Terphenyl	230 C18H14	000084-15-1 53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
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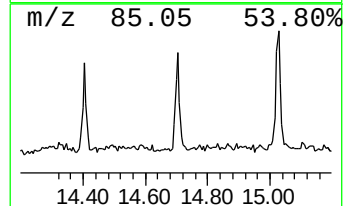
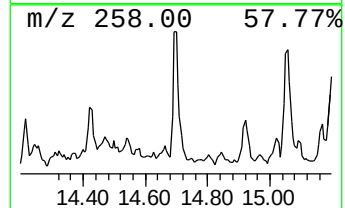
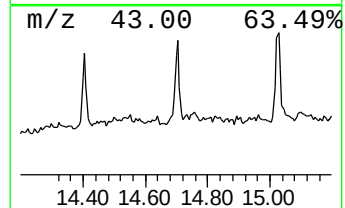
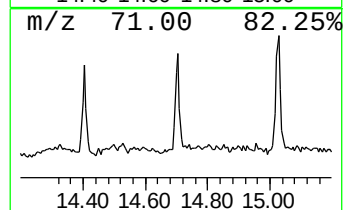
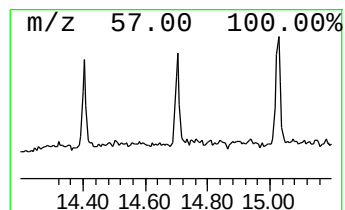
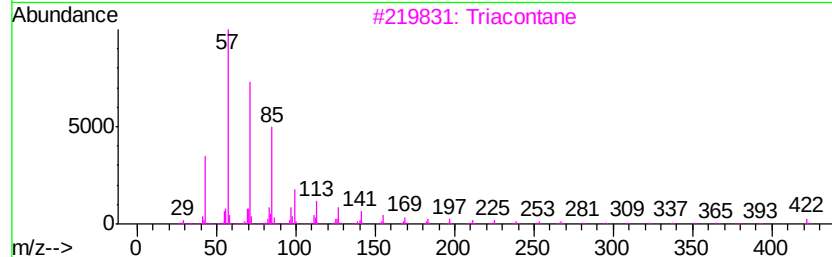
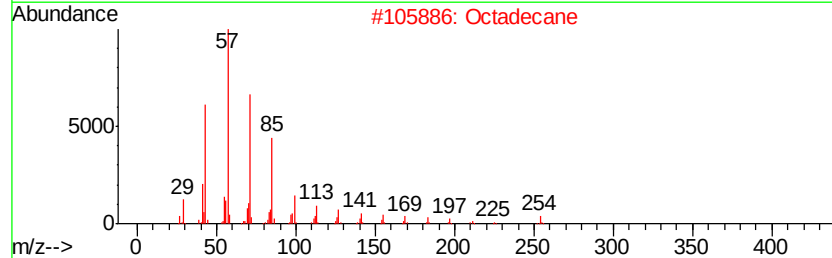
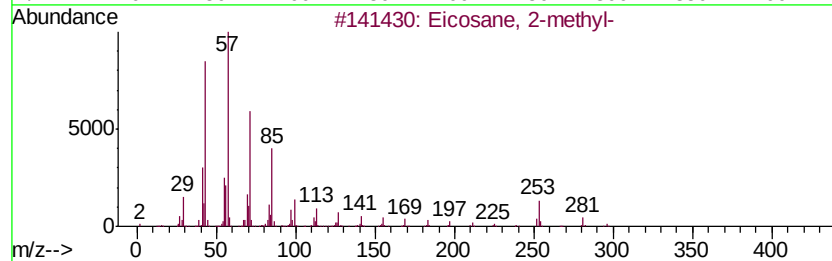
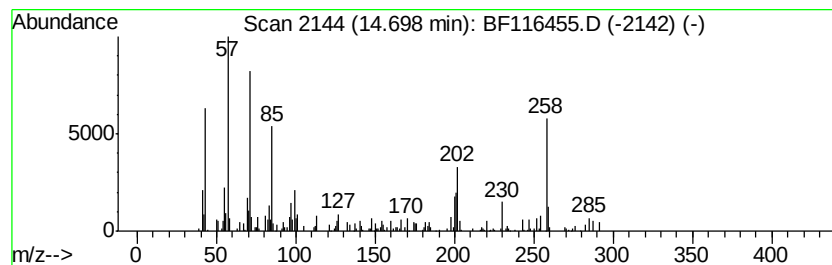
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	2.84 ng	101199	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane, 2-methyl-	296 C21H44	001560-84-5	68
2	Octadecane	254 C18H38	000593-45-3	64
3	Triacontane	422 C30H62	000638-68-6	62
4	Tetradecane	198 C14H30	000629-59-4	62
5	Octacosane	394 C28H58	000630-02-4	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
Data File : BF116455.D
Acq On : 21 Aug 2019 2:42
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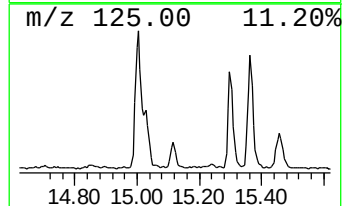
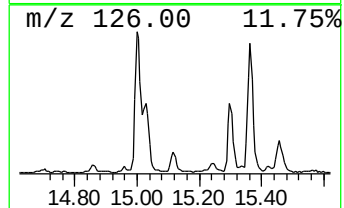
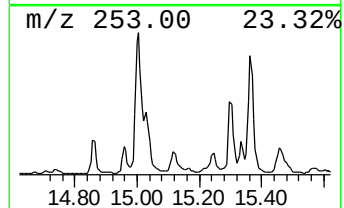
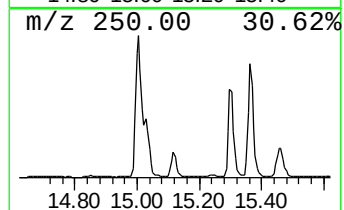
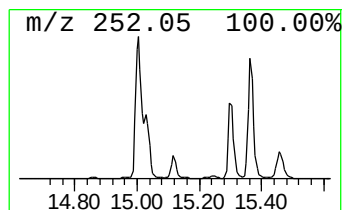
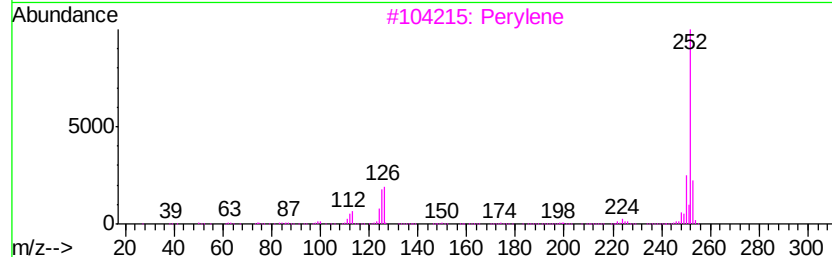
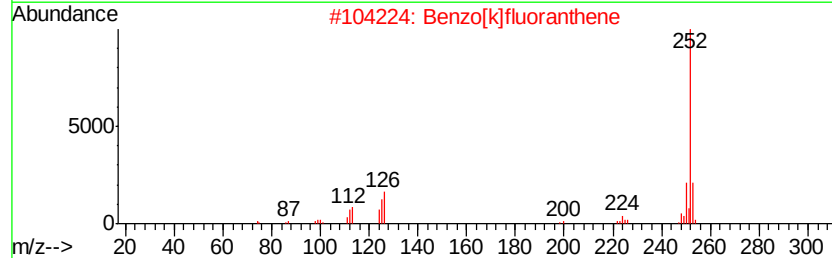
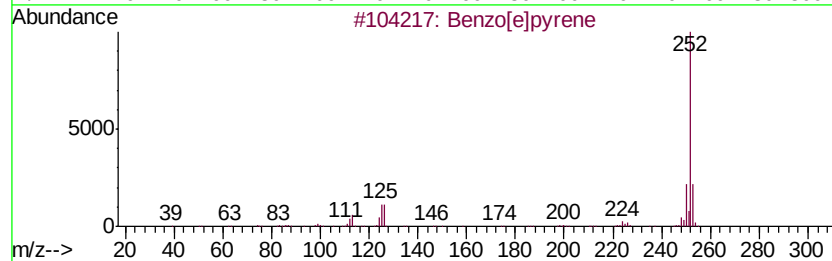
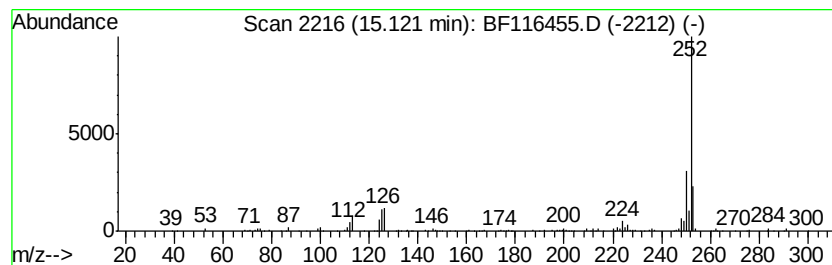
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	3.80 ng	135643	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3		Perylene	252	C20H12	000198-55-0	95
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[a]pyrene	252	C20H12	000050-32-8	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
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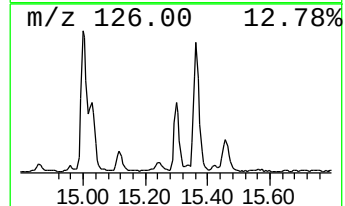
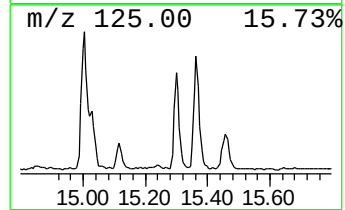
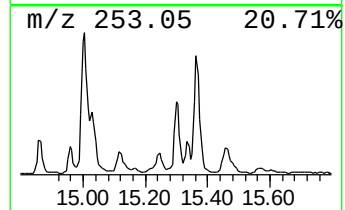
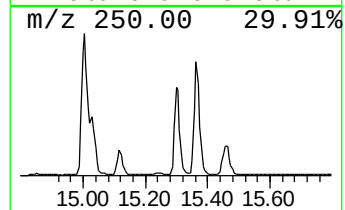
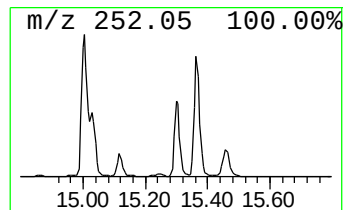
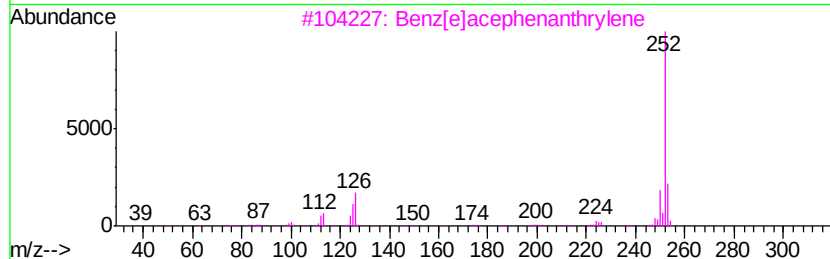
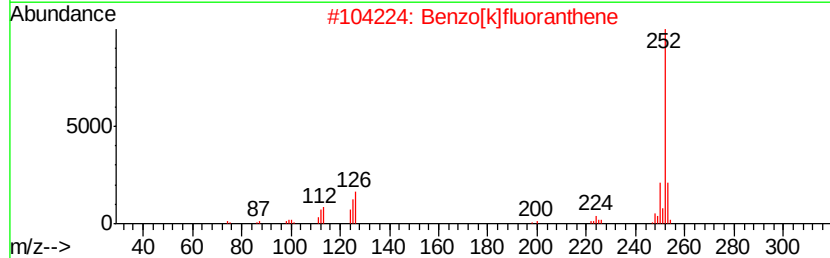
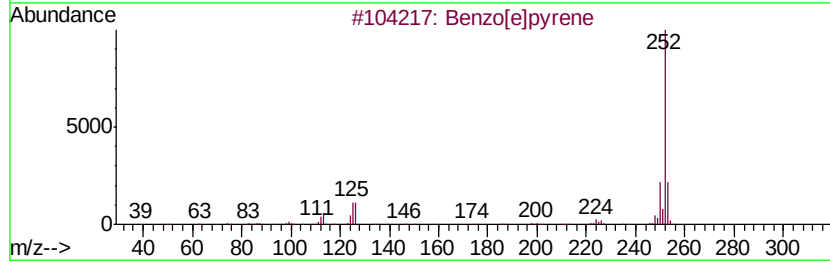
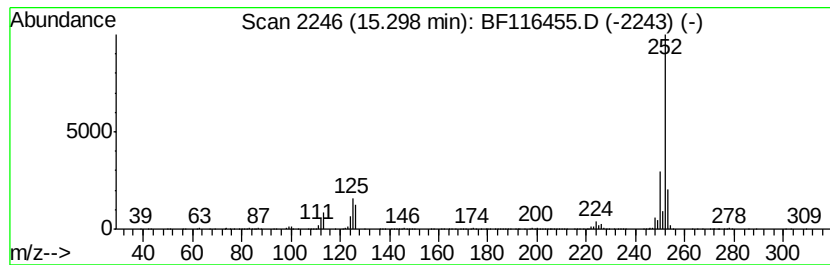
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Perylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.30	11.61 ng	414304	Perylene-d12	15.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[el]pyrene	252	C20H12	000192-97-2	99
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3	Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	98
4	Perylene	252	C20H12	000198-55-0	96
5	Benzo[a]pyrene	252	C20H12	000050-32-8	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082019\
Data File : BF116455.D
Acq On : 21 Aug 2019 2:42
Operator : HP/JU
Sample : K4354-02 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PX-02-081319

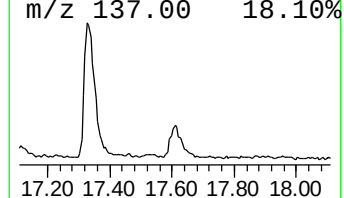
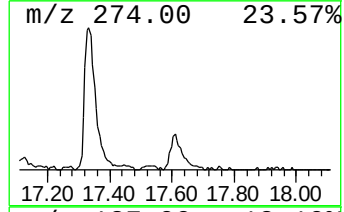
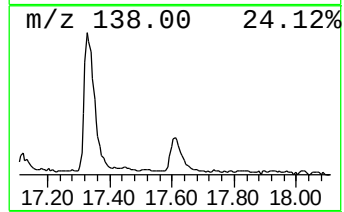
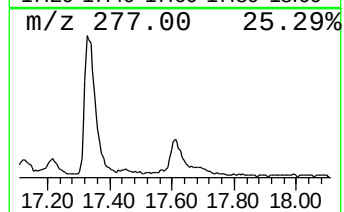
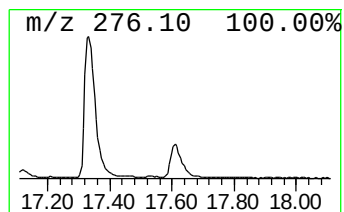
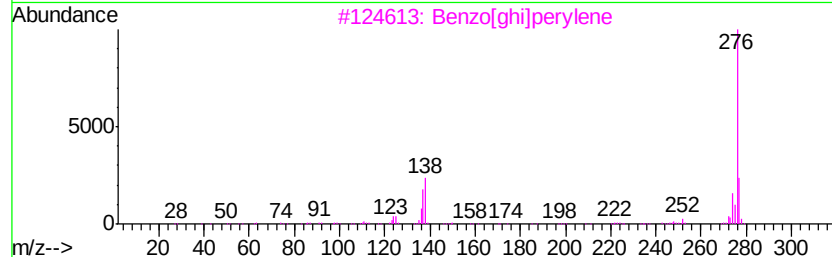
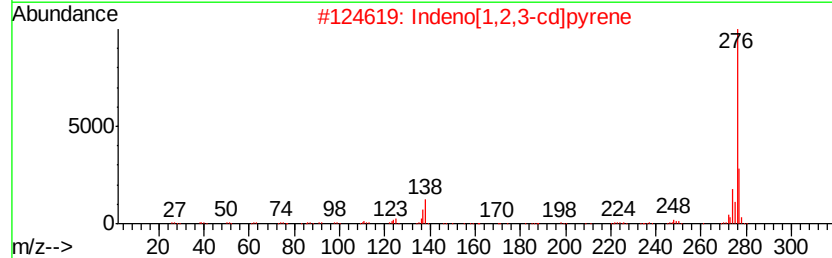
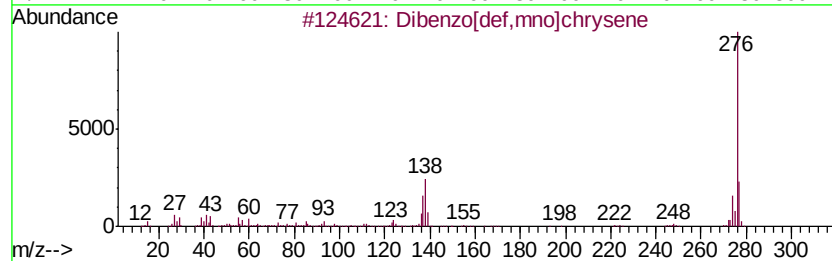
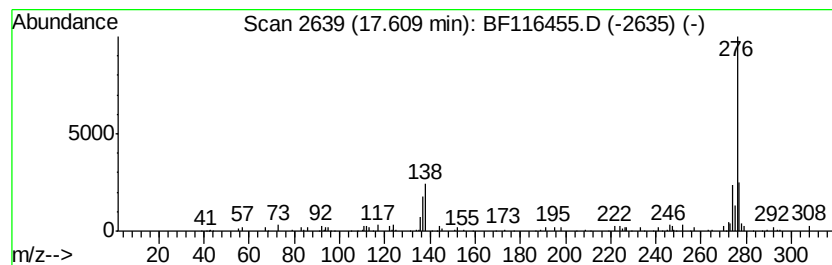
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dibenzo[def,mno]chrysene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.61	2.89 ng	103064	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	94
2		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	93
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	87
4		Naphthalen-1-yl-(3-nitro-benzyl)...	276	C17H12N2O2	200421-53-0	50
5		2-Amino-3-cyano-4,5-diphenylthio...	276	C17H12N2S	042160-34-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082019\
 Data File : BF116455.D
 Acq On : 21 Aug 2019 2:42
 Operator : HP/JU
 Sample : K4354-02 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PX-02-081319

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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
1,3,5-Cycloheptat...	3.81	8.1 ng		222740	1	6.80	549384	20.0
unknown6.58	6.58	38.4 ng		1053870	1	6.80	549384	20.0
Phenanthrene, 2-m...	11.83	2.0 ng		108853	4	11.32	1086740	20.0
Phenanthrene, 1-m...	11.86	3.2 ng		171752	4	11.32	1086740	20.0
4H-Cyclopenta[def...	11.94	4.2 ng		225987	4	11.32	1086740	20.0
11H-Benzo[b]fluorene	13.10	2.1 ng		210348	5	13.96	1978010	20.0
Pyrene, 1-methyl-	13.20	2.3 ng		228893	5	13.96	1978010	20.0
7H-Benz[de]anthra...	13.83	2.1 ng		205216	5	13.96	1978010	20.0
Eicosane, 2-methyl-	14.70	2.8 ng		101199	6	15.42	713746	20.0
Benzo[e]pyrene	15.12	3.8 ng		135643	6	15.42	713746	20.0
Perylene	15.30	11.6 ng		414304	6	15.42	713746	20.0
Dibenzo[def,mno]c...	17.61	2.9 ng		103064	6	15.42	713746	20.0