

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
 Data File : BF139442.D
 Acq On : 18 Sep 2024 23:07
 Operator : RC/JU
 Sample : P4086-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SOIL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139442.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.199	9	18	24	rBV	706606	1131881	69.96%	7.575%
2	3.975	313	320	326	rVB	15510	24011	1.48%	0.161%
3	5.099	505	511	518	rVB	99668	148490	9.18%	0.994%
4	5.499	569	579	584	rBV	1194628	1568791	96.96%	10.499%
5	6.504	744	750	756	rBV	1176434	1608305	99.40%	10.764%
6	6.875	808	813	820	rVB	319797	401834	24.84%	2.689%
7	7.440	898	909	914	rBV	825179	1073679	66.36%	7.186%
8	7.693	948	952	956	rVB	26069	32831	2.03%	0.220%
9	8.157	1024	1031	1038	rBV	374900	485601	30.01%	3.250%
10	8.575	1098	1102	1112	rVB	12663	21573	1.33%	0.144%
11	9.234	1208	1214	1219	rBV	1320688	1617973	100.00%	10.829%
12	9.310	1223	1227	1231	rBV	16025	20698	1.28%	0.139%
13	9.628	1277	1281	1288	rVB3	9774	17681	1.09%	0.118%
14	9.839	1312	1317	1321	rVB	19593	24450	1.51%	0.164%
15	9.910	1321	1329	1333	rBV	348544	427158	26.40%	2.859%
16	10.004	1338	1345	1351	rBV	138808	208613	12.89%	1.396%
17	10.334	1394	1401	1405	rBV2	17259	30132	1.86%	0.202%
18	10.622	1445	1450	1457	rVB	110854	137016	8.47%	0.917%
19	10.698	1457	1463	1468	rBV	615533	775962	47.96%	5.193%
20	10.816	1478	1483	1491	rVB2	27363	46934	2.90%	0.314%
21	11.257	1552	1558	1561	rBV2	19466	28608	1.77%	0.191%
22	11.398	1577	1582	1585	rBV	293502	448952	27.75%	3.005%
23	11.469	1590	1594	1597	rVV	27055	33363	2.06%	0.223%
24	11.628	1617	1621	1627	rBV3	11081	21113	1.30%	0.141%
25	11.686	1627	1631	1635	rVB2	18283	22792	1.41%	0.153%
26	11.922	1663	1671	1676	rBV2	205535	314938	19.46%	2.108%
27	12.004	1682	1685	1693	rVB2	45933	92393	5.71%	0.618%
28	12.092	1696	1700	1704	rBV	20237	26816	1.66%	0.179%
29	12.192	1713	1717	1719	rBV	20956	27695	1.71%	0.185%
30	12.339	1736	1742	1745	rBV2	17581	31956	1.98%	0.214%
31	12.445	1756	1760	1763	rBV	34604	46898	2.90%	0.314%
32	12.480	1763	1766	1772	rVV2	29147	44495	2.75%	0.298%
33	12.527	1772	1774	1780	rVB3	18691	26891	1.66%	0.180%
34	12.610	1783	1788	1792	rBV	252716	329332	20.35%	2.204%
35	12.704	1800	1804	1811	rBV2	52418	86970	5.38%	0.582%
36	12.839	1820	1827	1832	rVV	229297	354452	21.91%	2.372%

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ALS Vial : 16 Sample Multiplier: 1

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.886	1832	1835	1840	rVB2	21894	33911	2.10%	0.227%
38	12.980	1843	1851	1858	rVB	951197	1217332	75.24%	8.147%
39	13.051	1859	1863	1870	rBV4	28810	56079	3.47%	0.375%
40	13.163	1879	1882	1886	rVB	46170	60493	3.74%	0.405%
41	13.227	1891	1893	1896	rVV	18529	19735	1.22%	0.132%
42	13.263	1896	1899	1903	rVB2	33652	42783	2.64%	0.286%
43	13.392	1918	1921	1924	rBV2	17281	20198	1.25%	0.135%
44	13.669	1965	1968	1976	rVB	35098	50753	3.14%	0.340%
45	13.780	1983	1987	1990	rBV2	29354	47601	2.94%	0.319%
46	13.874	1999	2003	2011	rVB2	93095	156793	9.69%	1.049%
47	14.027	2023	2029	2033	rBV2	329298	586072	36.22%	3.922%
48	15.074	2202	2207	2215	rVB	111805	254017	15.70%	1.700%
49	15.374	2254	2258	2264	rVV	60994	94420	5.84%	0.632%
50	15.433	2264	2268	2274	rVB	87539	135074	8.35%	0.904%
51	15.504	2275	2280	2293	rVB2	140197	275957	17.06%	1.847%
52	16.962	2523	2528	2537	rVB2	38785	85948	5.31%	0.575%
53	17.410	2598	2604	2613	rVB	27140	63317	3.91%	0.424%

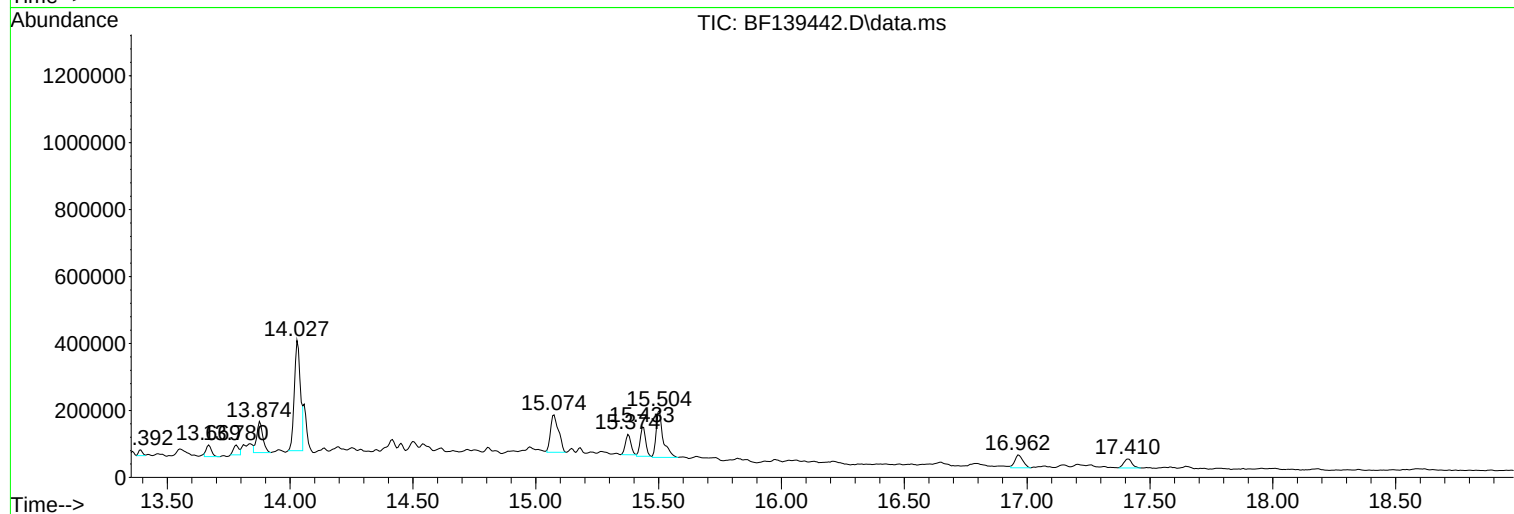
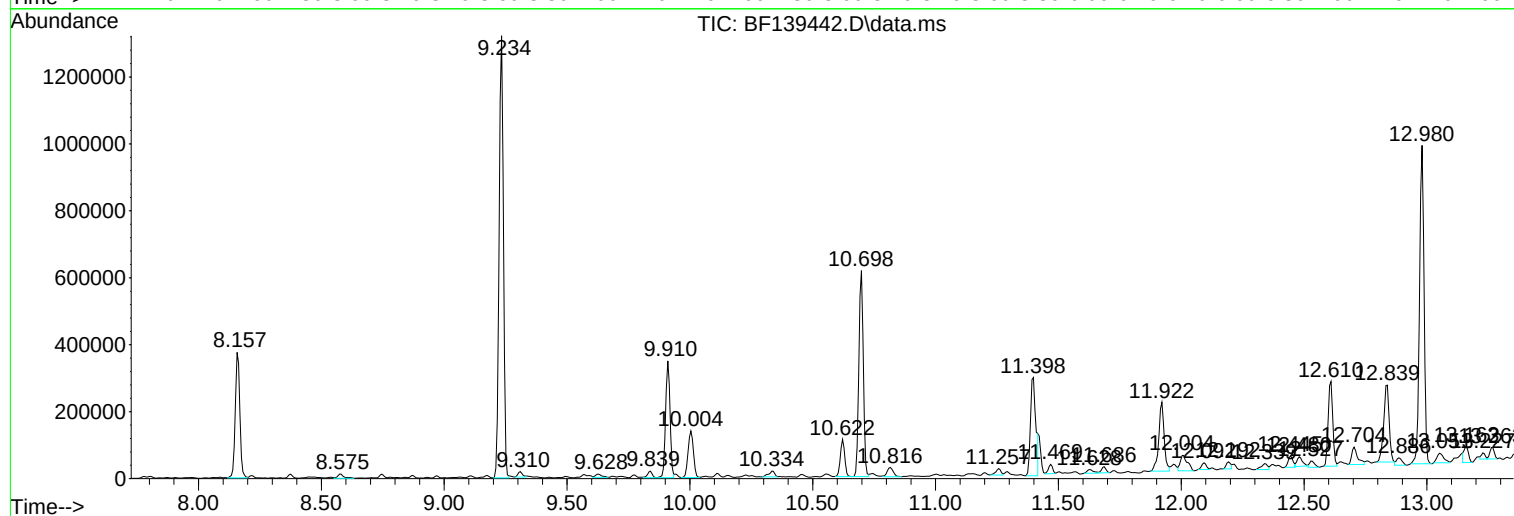
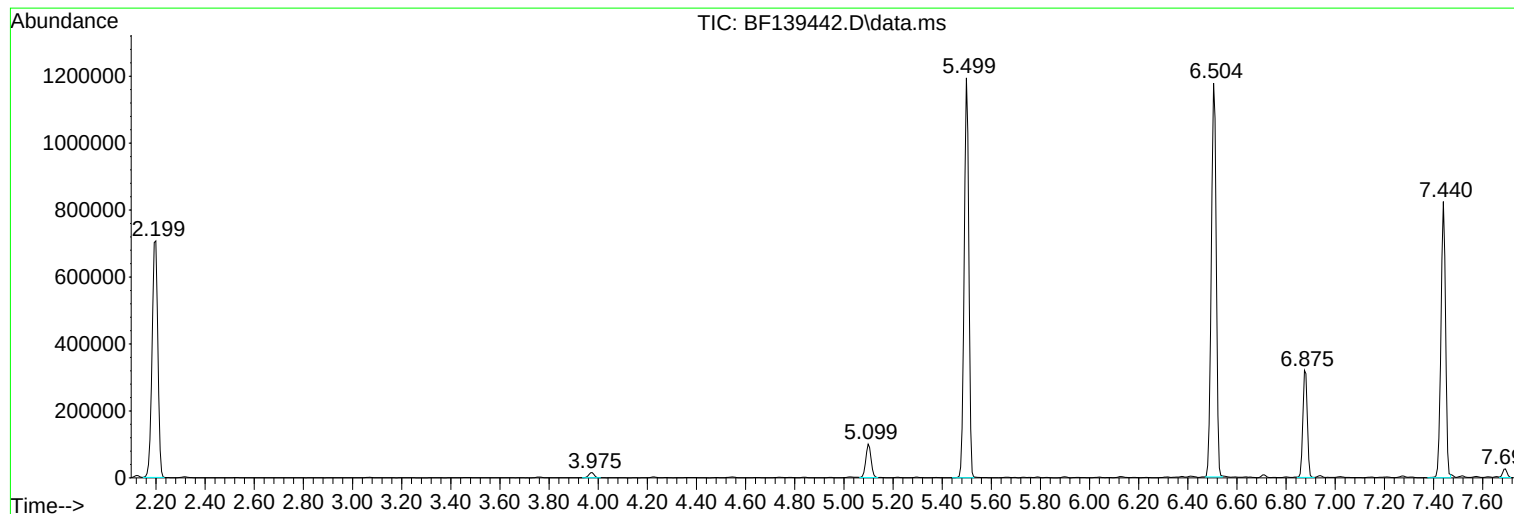
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
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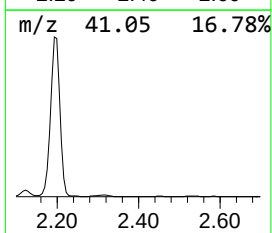
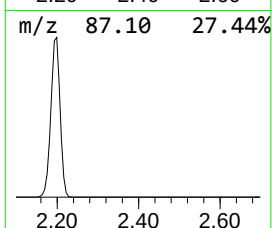
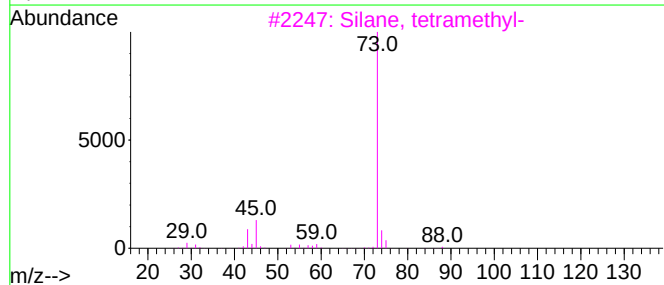
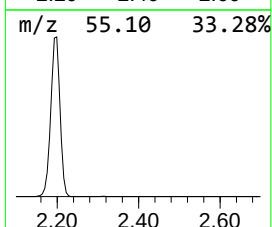
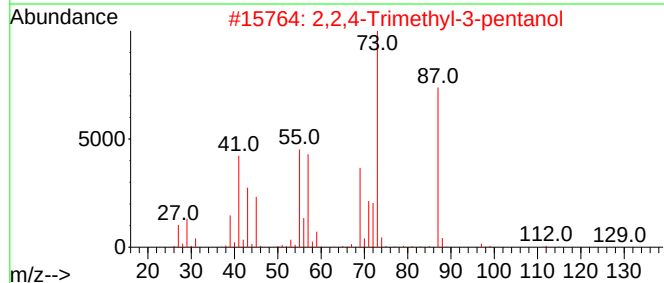
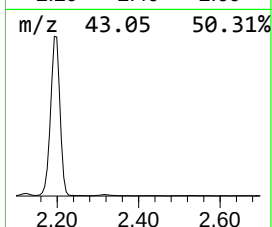
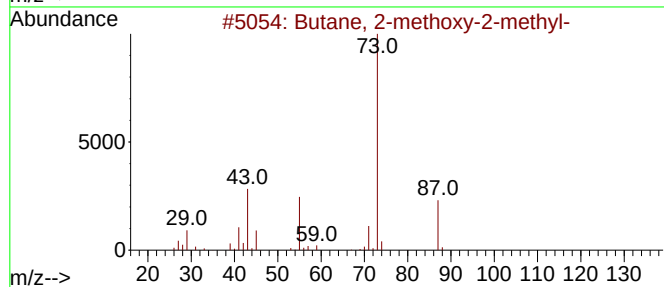
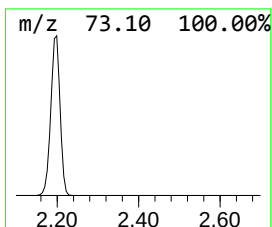
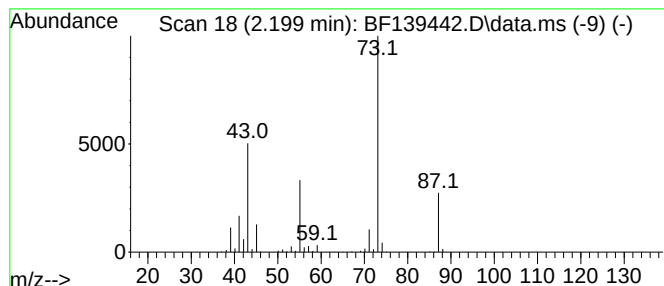
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.199	56.34 ng	1131880	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			2-OXOPROPIONAMIDE	87	C3H5NO2	000631-66-3	12
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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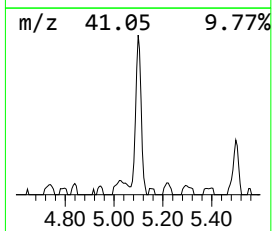
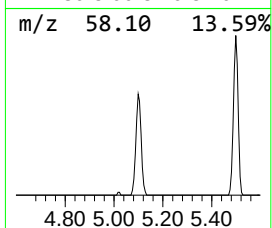
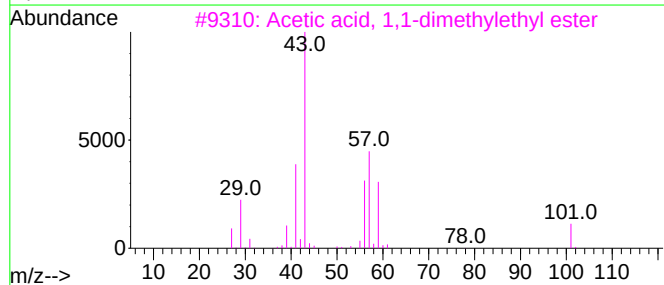
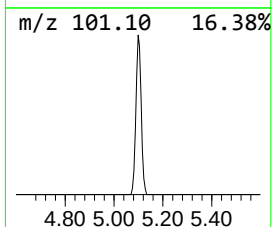
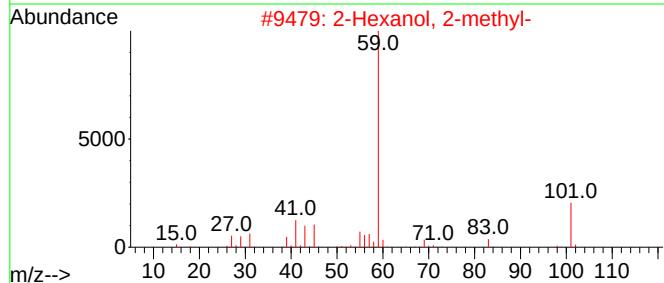
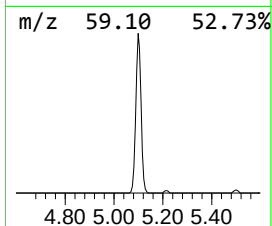
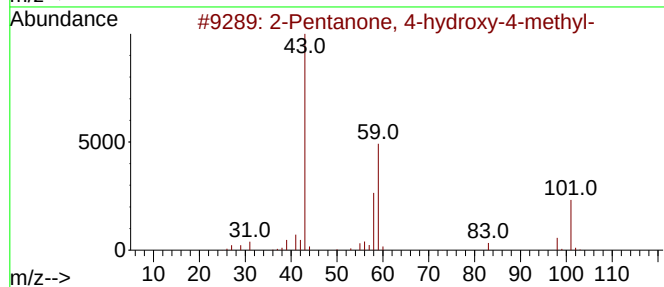
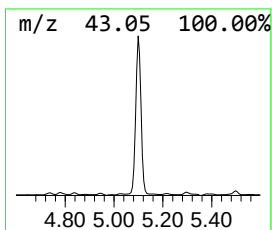
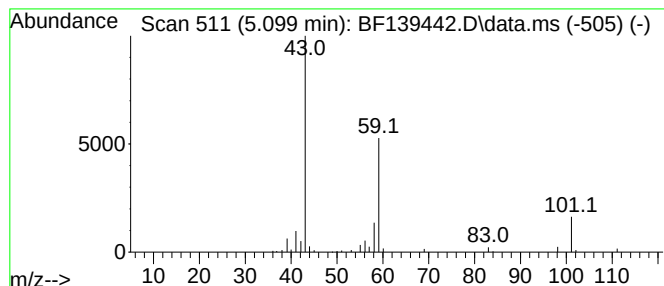
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.099	7.39 ng	148490	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	37		
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23		



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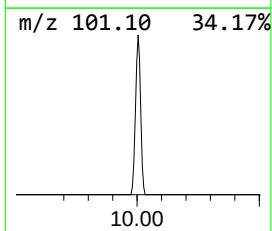
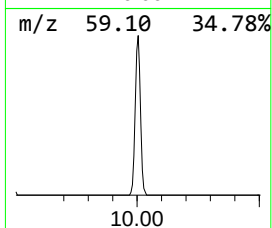
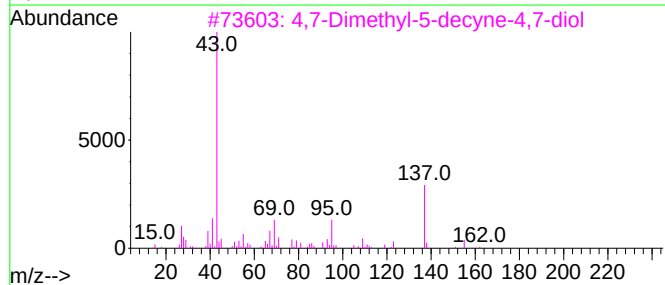
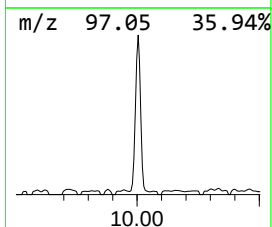
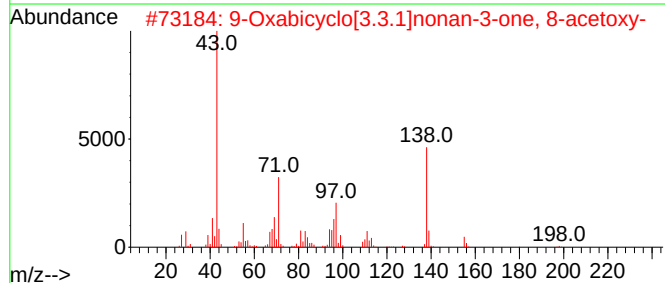
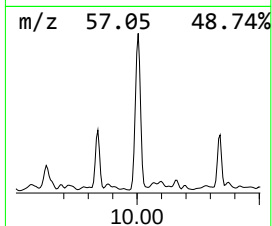
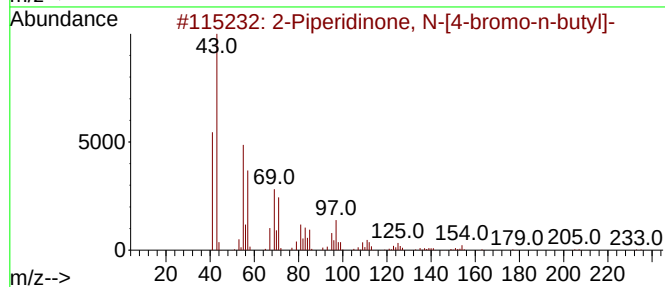
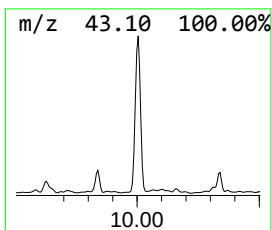
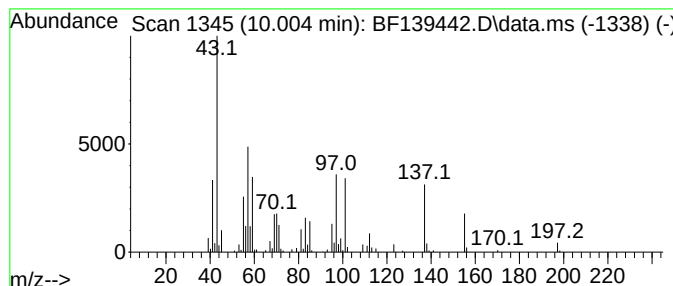
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown10.004 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.004	9.77 ng	208613	Acenaphthene-d10	9.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	14		
2	9-Oxabicyclo[3.3.1]nonan-3-one, ...	198	C10H14O4	1010197-59-0	14		
3	4,7-Dimethyl-5-decyne-4,7-diol	198	C12H22O2	000126-87-4	12		
4	2-Pentanol, 2-methyl-, acetate	144	C8H16O2	034859-98-8	10		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10		



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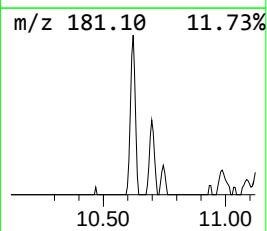
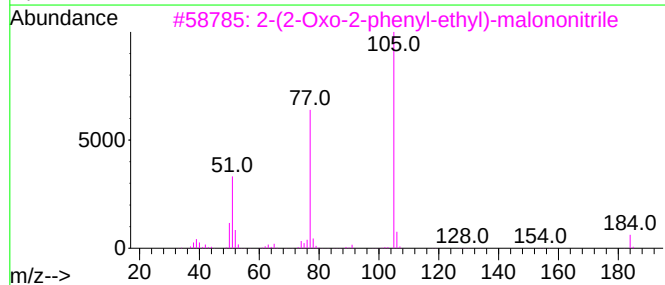
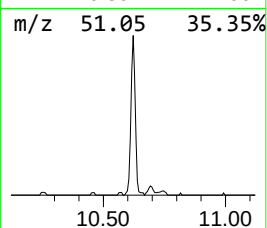
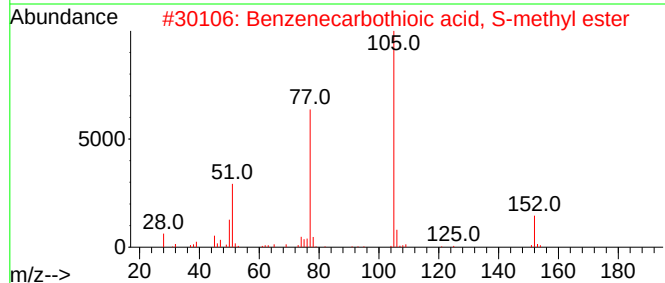
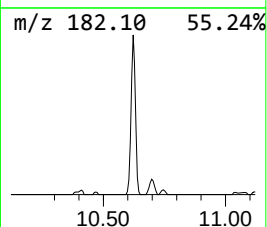
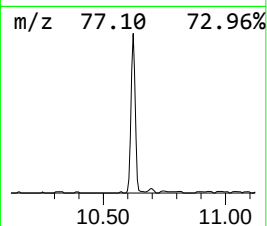
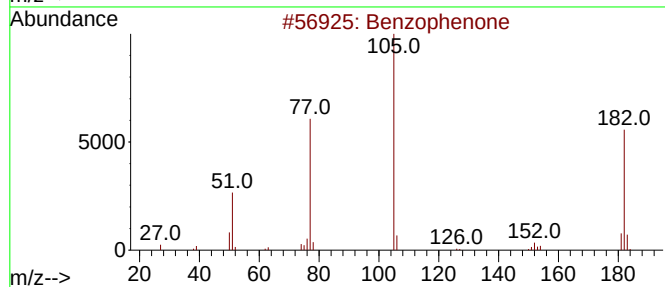
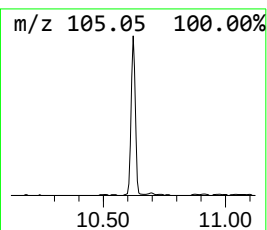
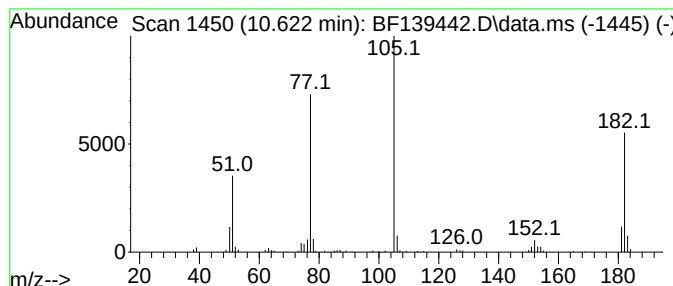
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.622	6.42 ng	137016	Acenaphthene-d10	9.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	60
3			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47
4			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
5			Methyl 2-benzoyl-4-cyanobutanoate	231	C13H13NO3	1010486-83-8	47



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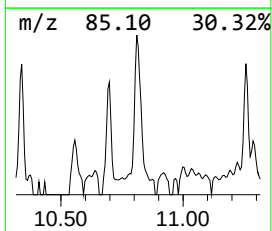
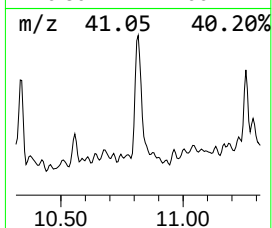
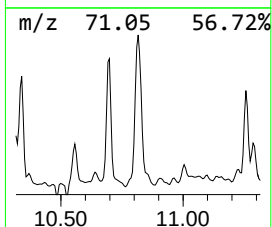
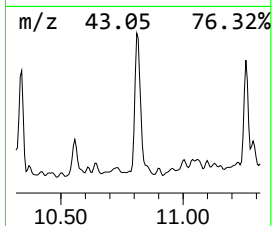
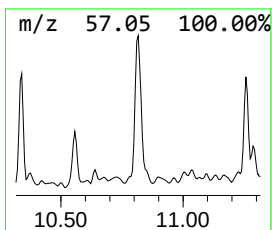
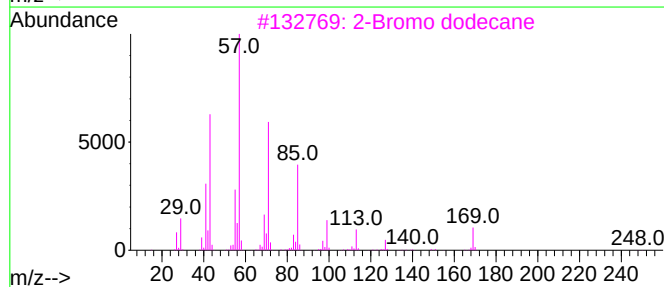
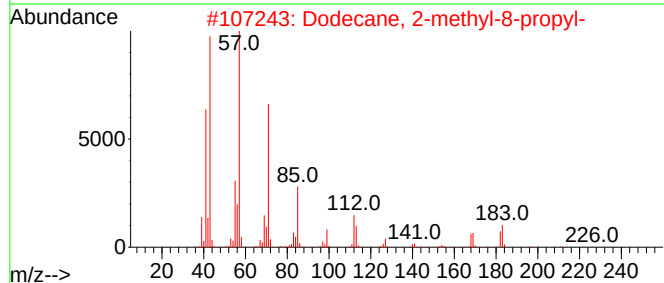
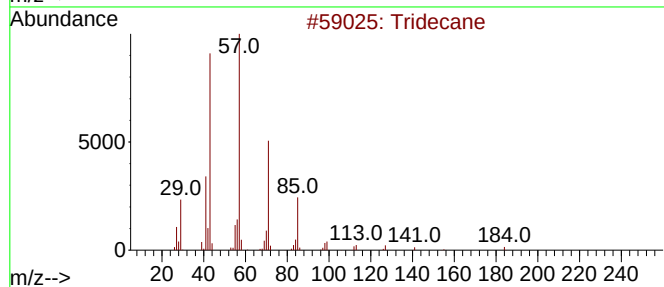
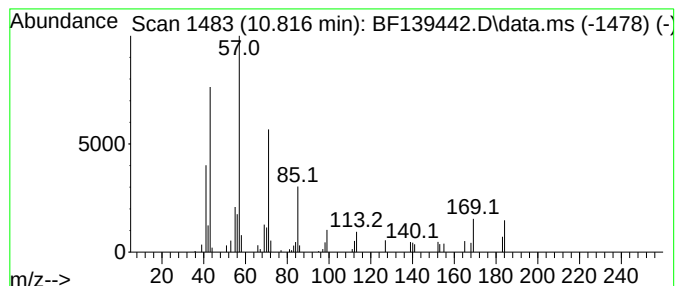
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	2.09 ng	46934	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	81
2			Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	80
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	72
4			Heneicosane	296	C21H44	000629-94-7	68
5			Tetratetracontane	619	C44H90	007098-22-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
Operator : RC/JU
Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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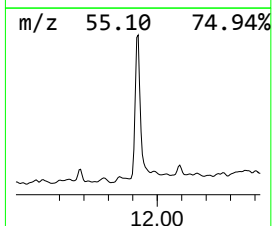
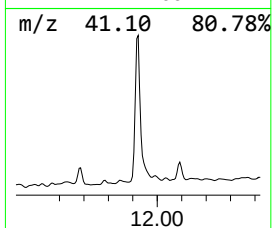
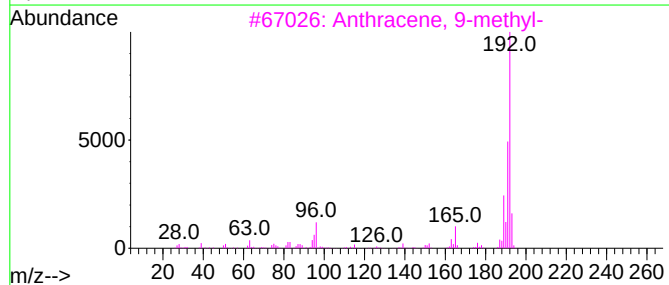
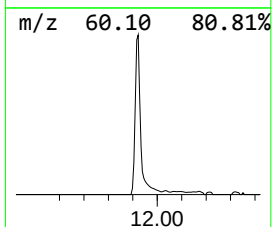
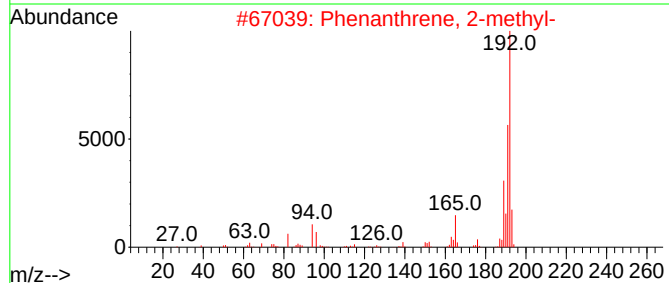
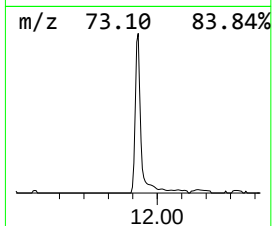
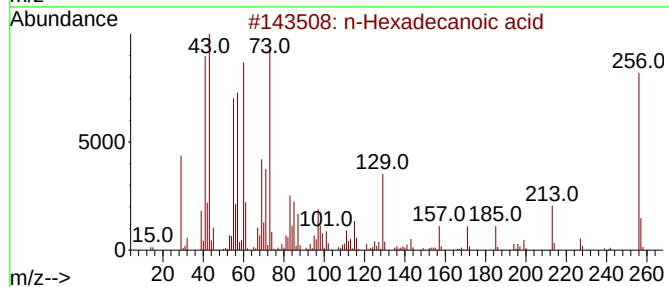
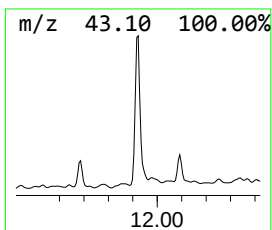
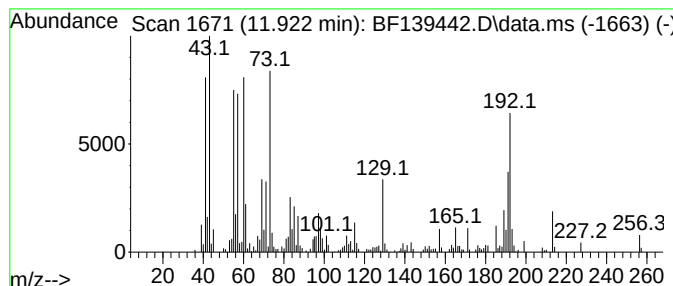
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.922	14.03 ng	314938	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
4			n-Decanoic acid	172	C10H20O2	000334-48-5	89
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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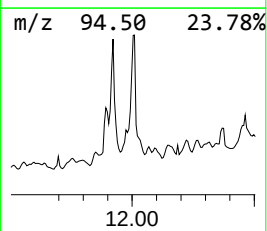
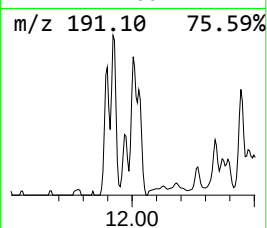
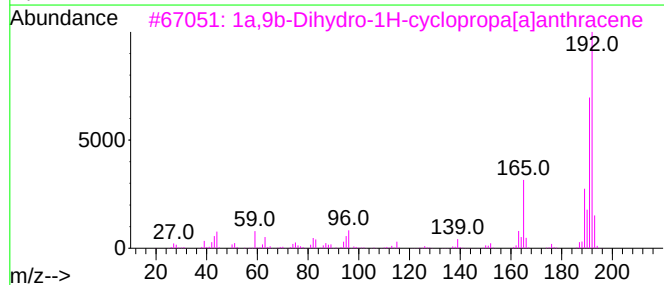
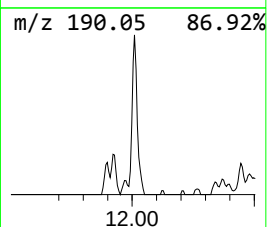
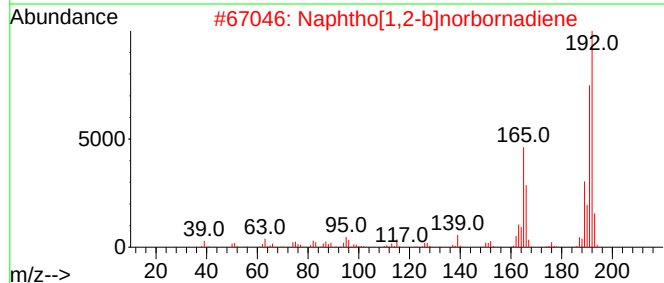
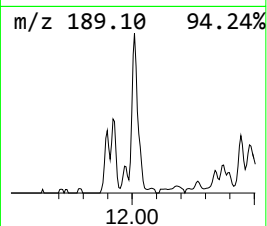
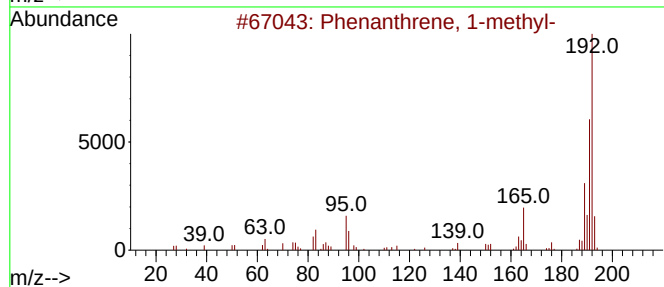
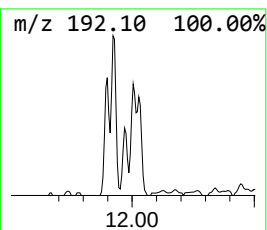
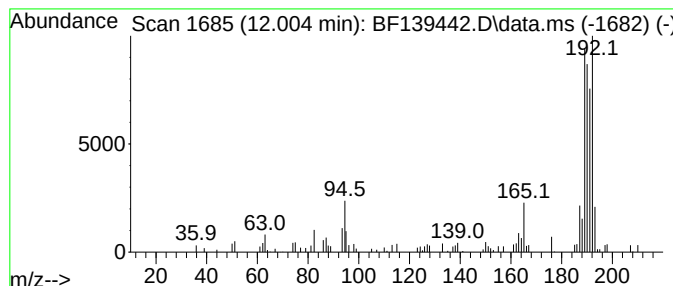
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	4.12 ng	92393	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	50
2			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	50
3			1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	46
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	45
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
Operator : RC/JU
Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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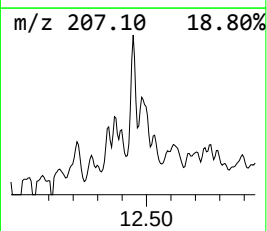
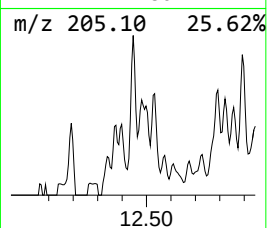
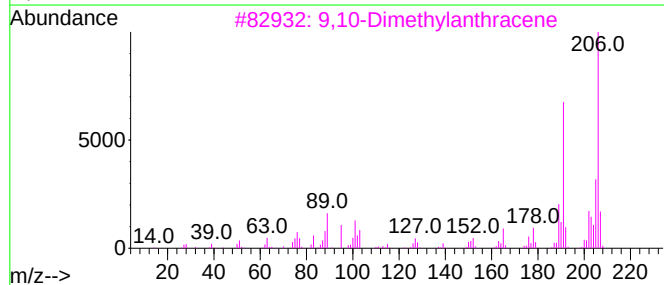
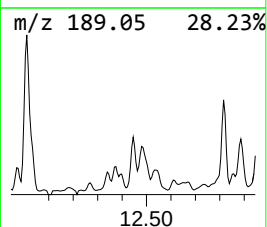
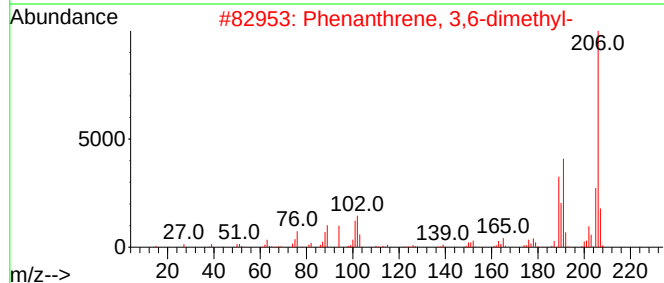
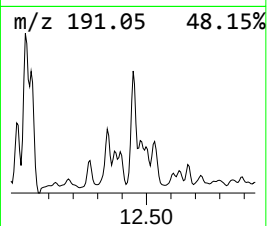
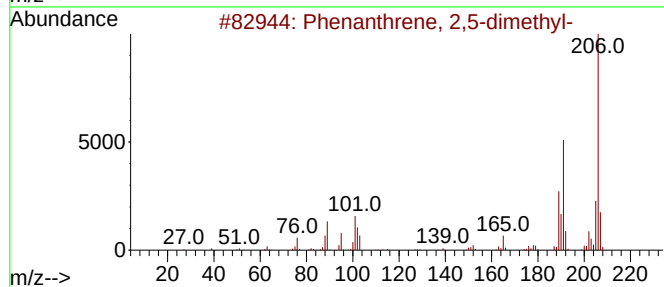
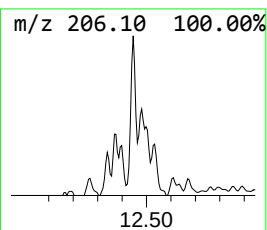
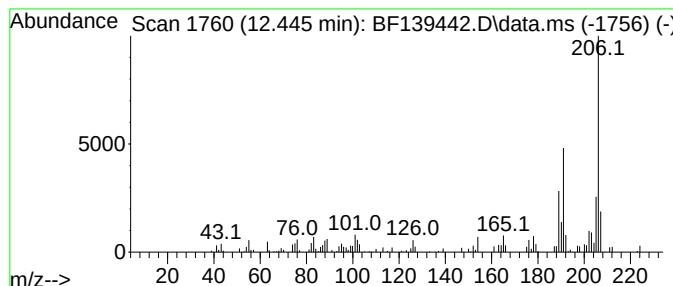
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	2.09 ng	46898	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3			9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
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ALS Vial : 16 Sample Multiplier: 1

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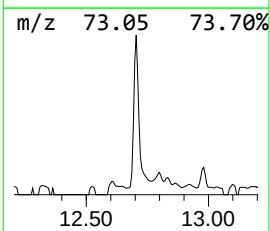
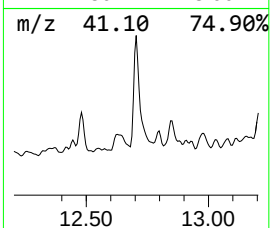
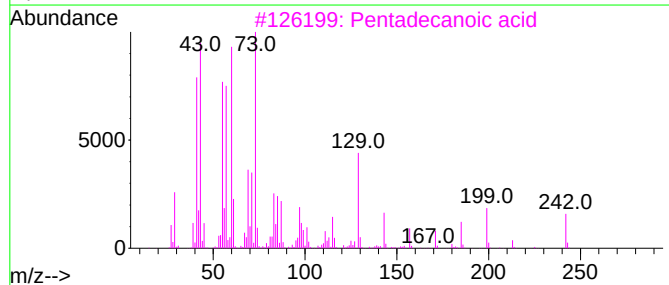
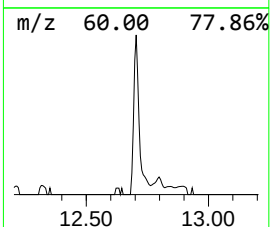
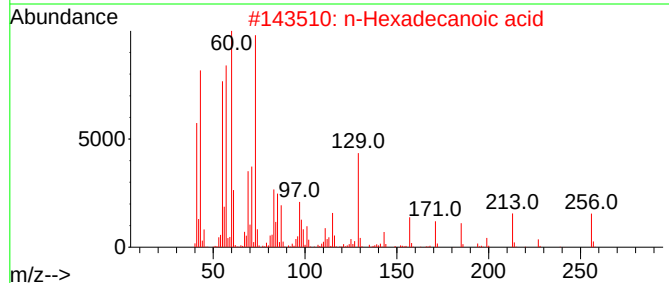
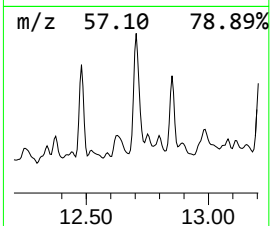
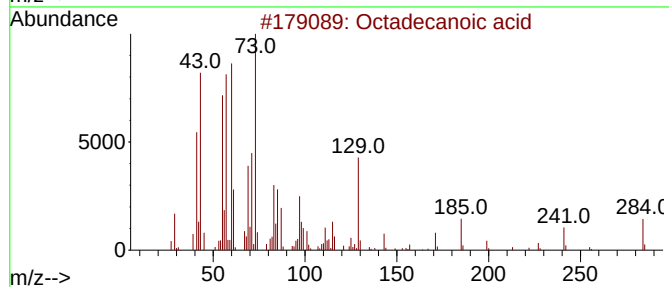
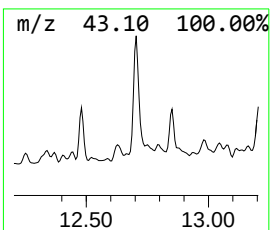
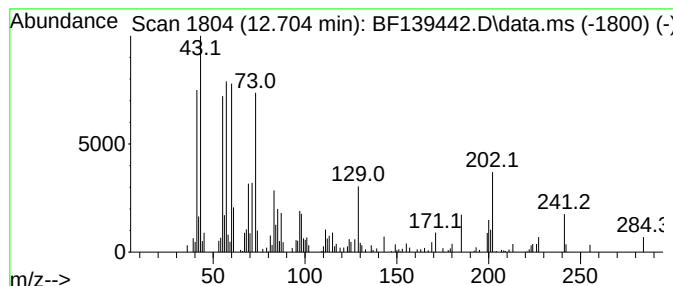
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.704	3.87 ng	86970	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	91
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4			Dodecanoic acid	200	C12H24O2	000143-07-7	64
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
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Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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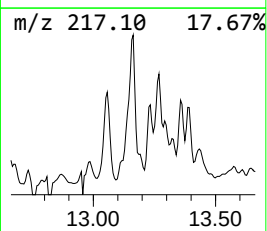
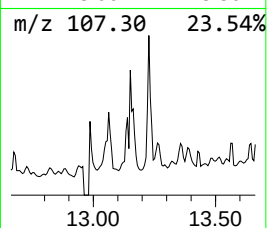
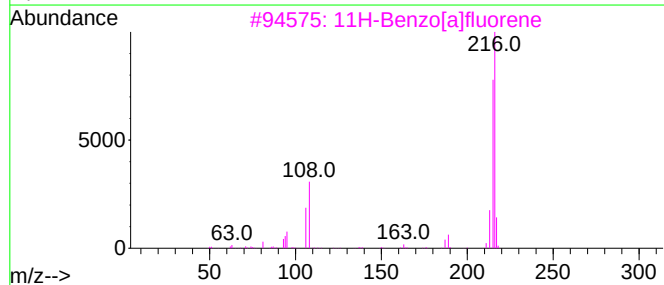
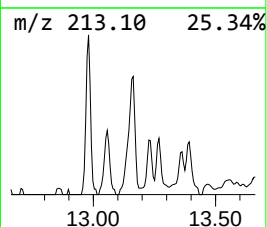
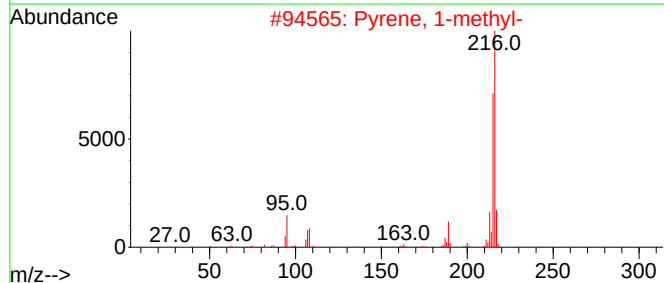
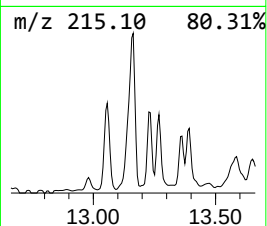
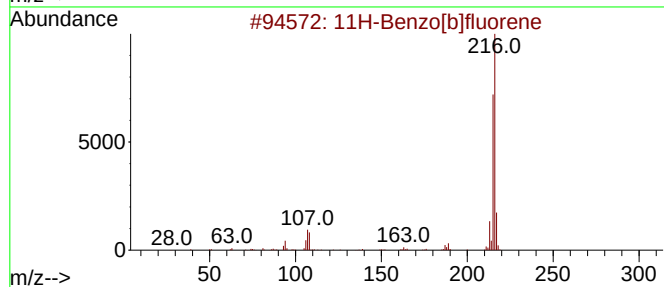
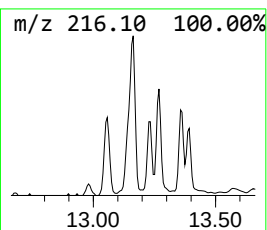
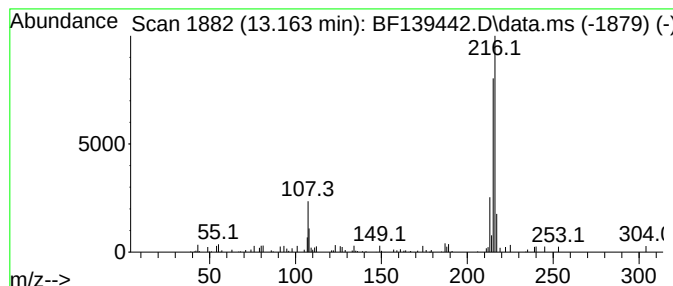
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	2.06 ng	60493	Chrysene-d12	14.033

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64
5		1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
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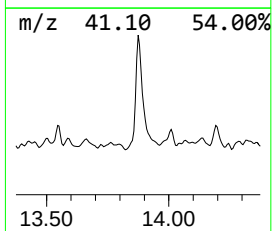
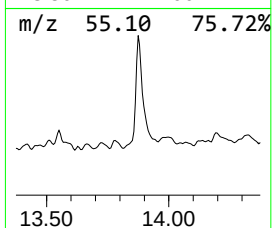
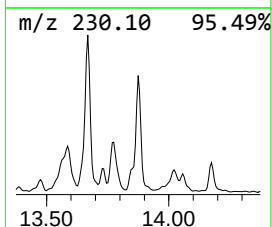
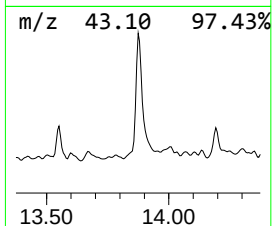
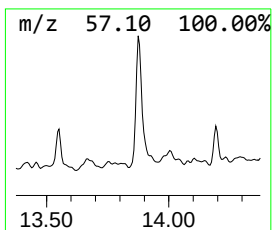
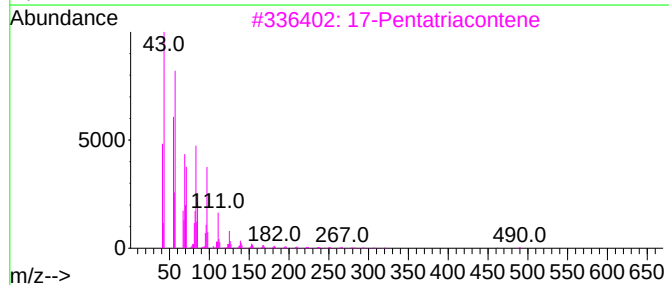
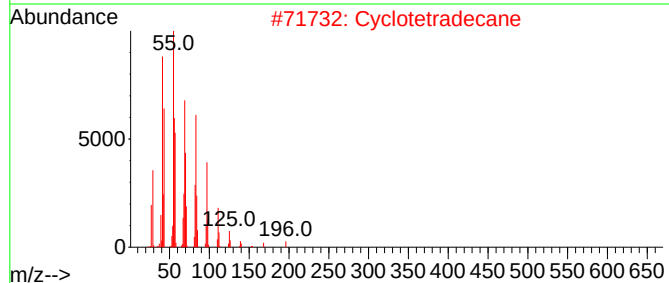
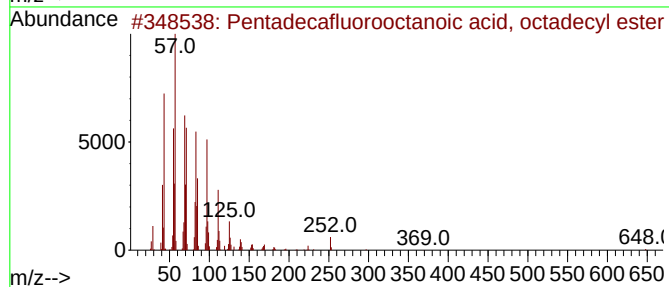
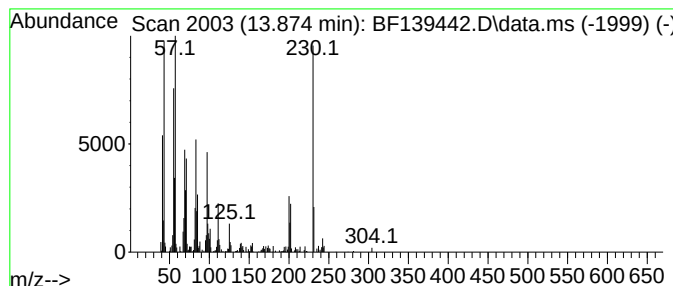
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	5.35 ng	156793	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	87
2			Cyclotetradecane	196	C14H28	000295-17-0	59
3			17-Pentatriacontene	491	C35H70	006971-40-0	50
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	50
5			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
Operator : RC/JU
Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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SOIL

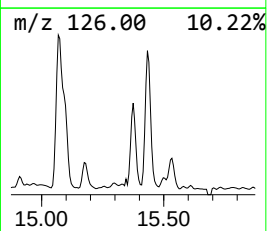
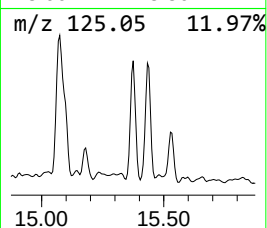
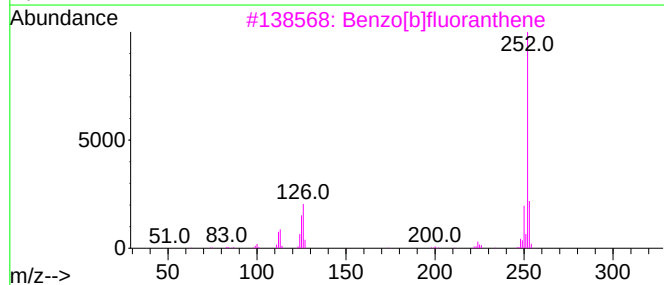
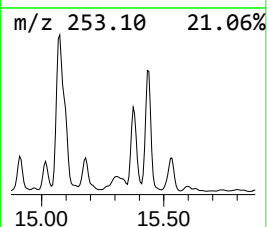
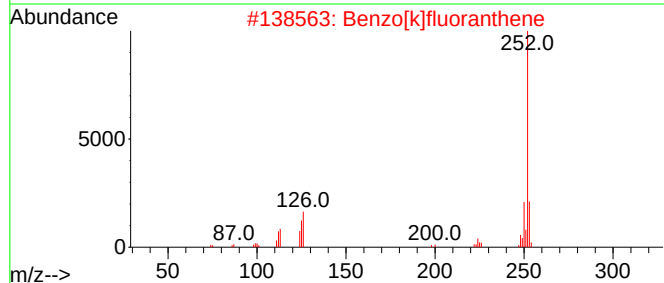
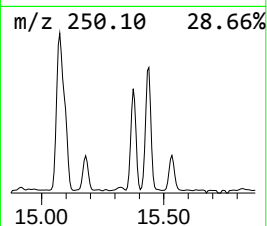
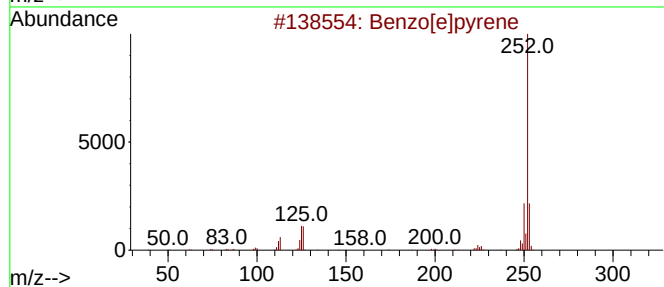
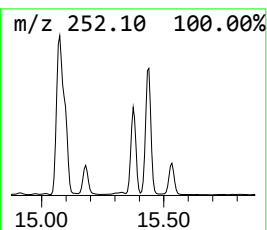
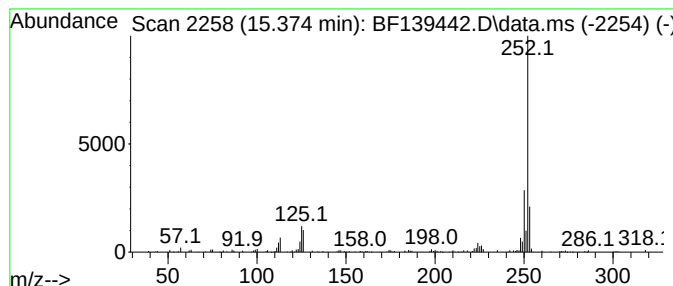
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.374	6.84 ng	94420	Perylene-d12	15.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	97
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90
5			Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091824\
Data File : BF139442.D
Acq On : 18 Sep 2024 23:07
Operator : RC/JU
Sample : P4086-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SOIL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.199	56.3	ng	1131880	1	6.875	401834	20.0
2-Pentanone, 4-...	5.099	7.4	ng	148490	1	6.875	401834	20.0
unknown10.004	10.004	9.8	ng	208613	3	9.910	427158	20.0
Benzophenone	10.622	6.4	ng	137016	3	9.910	427158	20.0
Tridecane	10.816	2.1	ng	46934	4	11.398	448952	20.0
n-Hexadecanoic ...	11.922	14.0	ng	314938	4	11.398	448952	20.0
Phenanthrene, 1...	12.004	4.1	ng	92393	4	11.398	448952	20.0
Phenanthrene, 2...	12.445	2.1	ng	46898	4	11.398	448952	20.0
Octadecanoic acid	12.704	3.9	ng	86970	4	11.398	448952	20.0
11H-Benzo[b]flu...	13.163	2.1	ng	60493	5	14.033	586072	20.0
Pentadecafluoro...	13.874	5.3	ng	156793	5	14.033	586072	20.0
Benzo[e]pyrene	15.374	6.8	ng	94420	6	15.504	275957	20.0