

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
 Data File : BF143707.D
 Acq On : 11 Sep 2025 15:32
 Operator : RC/JU
 Sample : Q3055-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-090925

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-BF091025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.110	487	496	501	rBB2	32081	49421	8.47%	0.618%
2	5.504	557	563	568	rBV	405250	520415	89.21%	6.506%
3	6.504	727	733	738	rBV	373005	495563	84.95%	6.195%
4	6.892	794	799	804	rVB	179867	213572	36.61%	2.670%
5	7.451	888	894	899	rBV	256918	324428	55.61%	4.056%
6	8.175	1012	1017	1025	rBB	229178	291584	49.98%	3.645%
7	8.886	1134	1138	1142	rBV	9510	11149	1.91%	0.139%
8	8.986	1150	1155	1160	rVB	11097	13324	2.28%	0.167%
9	9.245	1194	1199	1204	rBV	458726	583362	100.00%	7.293%
10	9.516	1239	1245	1248	rBV2	5427	7716	1.32%	0.096%
11	9.586	1253	1257	1260	rBV	10645	14101	2.42%	0.176%
12	9.704	1273	1277	1288	rVB4	4627	7857	1.35%	0.098%
13	9.792	1288	1292	1296	rBV	13437	18033	3.09%	0.225%
14	9.928	1308	1315	1319	rBV	243174	312670	53.60%	3.909%
15	9.963	1319	1321	1325	rVB	27298	27342	4.69%	0.342%
16	10.133	1345	1350	1354	rBV2	15537	19616	3.36%	0.245%
17	10.228	1360	1366	1372	rBV3	6590	13939	2.39%	0.174%
18	10.339	1380	1385	1390	rBV4	4999	10039	1.72%	0.125%
19	10.475	1404	1408	1414	rVB	34544	43390	7.44%	0.542%
20	10.539	1414	1419	1421	rBV2	4506	6977	1.20%	0.087%
21	10.639	1430	1436	1442	rVB	63011	84071	14.41%	1.051%
22	10.716	1444	1449	1454	rBV	273031	337565	57.87%	4.220%
23	10.839	1465	1470	1477	rVB3	8043	11348	1.95%	0.142%
24	10.951	1485	1489	1493	rVB	9267	11633	1.99%	0.145%
25	11.022	1493	1501	1504	rBV4	7580	13623	2.34%	0.170%
26	11.051	1504	1506	1512	rVV5	5026	7198	1.23%	0.090%
27	11.175	1519	1527	1531	rBV6	3652	9323	1.60%	0.117%
28	11.222	1531	1535	1539	rVB	5414	6855	1.18%	0.086%
29	11.316	1547	1551	1556	rVB	14915	18612	3.19%	0.233%
30	11.416	1564	1568	1570	rBV	201890	264082	45.27%	3.301%
31	11.439	1570	1572	1577	rVV	228856	276327	47.37%	3.454%
32	11.492	1577	1581	1584	rVV	41930	53640	9.19%	0.671%
33	11.522	1584	1586	1590	rVV2	7793	10166	1.74%	0.127%
34	11.645	1601	1607	1614	rVV	17962	27878	4.78%	0.349%
35	11.716	1615	1619	1622	rVV2	7378	11202	1.92%	0.140%
36	11.745	1622	1624	1628	rVV2	7561	9439	1.62%	0.118%

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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-BF091025.M
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37	11.798	1628	1633	1636	rVV5	8447	11999	2.06%	0.150%
38	11.939	1648	1657	1663	rBV2	109658	191686	32.86%	2.396%
39	11.992	1663	1666	1669	rVV	10366	16111	2.76%	0.201%
40	12.033	1669	1673	1681	rVB	57060	101802	17.45%	1.273%
41	12.116	1682	1687	1690	rBV3	4715	8463	1.45%	0.106%
42	12.151	1690	1693	1696	rVV3	5294	7033	1.21%	0.088%
43	12.216	1699	1704	1714	rVB3	23658	51918	8.90%	0.649%
44	12.363	1723	1729	1732	rBV5	5891	11681	2.00%	0.146%
45	12.469	1742	1747	1750	rBV	24419	32530	5.58%	0.407%
46	12.504	1750	1753	1758	rVV4	13973	20471	3.51%	0.256%
47	12.551	1758	1761	1770	rVB4	13404	22036	3.78%	0.275%
48	12.633	1770	1775	1780	rBV	326142	414836	71.11%	5.186%
49	12.686	1780	1784	1787	rBV4	4478	7509	1.29%	0.094%
50	12.727	1787	1791	1796	rBV2	33694	45795	7.85%	0.572%
51	12.786	1797	1801	1806	rVB3	13479	20379	3.49%	0.255%
52	12.863	1806	1814	1819	rBV	300132	442402	75.84%	5.530%
53	12.910	1819	1822	1827	rVB2	12121	18061	3.10%	0.226%
54	13.004	1830	1838	1845	rVB	386639	512690	87.89%	6.409%
55	13.074	1845	1850	1855	rBV4	23325	46557	7.98%	0.582%
56	13.186	1858	1869	1873	rBV	41335	80820	13.85%	1.010%
57	13.251	1873	1880	1884	rBV	32437	52582	9.01%	0.657%
58	13.292	1884	1887	1890	rVV2	23482	26727	4.58%	0.334%
59	13.386	1899	1903	1905	rBV	13493	15783	2.71%	0.197%
60	13.416	1905	1908	1911	rVB	16122	17195	2.95%	0.215%
61	13.498	1918	1922	1926	rBV6	6319	11916	2.04%	0.149%
62	13.692	1952	1955	1960	rVB	20887	30067	5.15%	0.376%
63	13.810	1969	1975	1977	rBV3	30323	50050	8.58%	0.626%
64	13.904	1987	1991	2000	rVB4	37399	61665	10.57%	0.771%
65	14.057	2011	2017	2020	rBV2	306026	514409	88.18%	6.431%
66	14.163	2032	2035	2039	rVB	24357	27445	4.70%	0.343%
67	14.274	2052	2054	2058	rBV2	9172	12317	2.11%	0.154%
68	14.439	2079	2082	2086	rVB	20523	26778	4.59%	0.335%
69	15.104	2190	2195	2205	rBV	125375	288904	49.52%	3.612%
70	15.410	2243	2247	2253	rVB	66403	107736	18.47%	1.347%
71	15.474	2253	2258	2264	rBV	91831	142840	24.49%	1.786%
72	15.539	2264	2269	2273	rBV	134334	207623	35.59%	2.595%
73	16.268	2390	2393	2402	rVB7	15594	32080	5.50%	0.401%
74	17.027	2517	2522	2533	rVB2	39674	93675	16.06%	1.171%
75	17.480	2594	2599	2613	rVB	32009	77318	13.25%	0.967%

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ClientSampleId :
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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-BF091025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

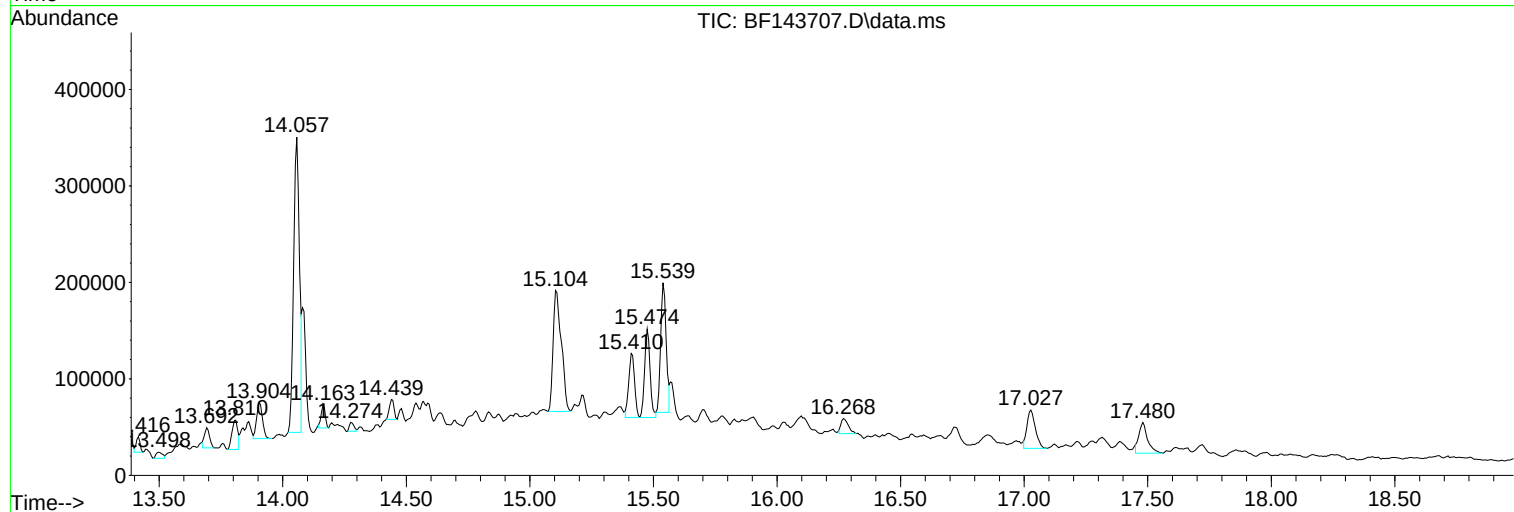
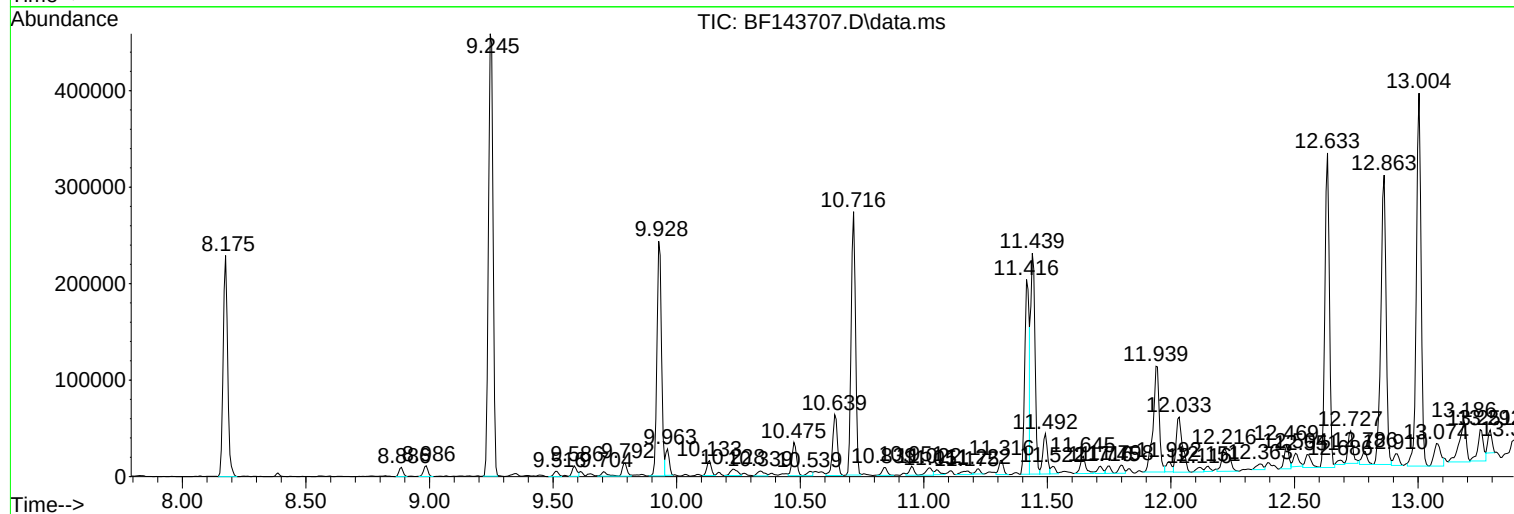
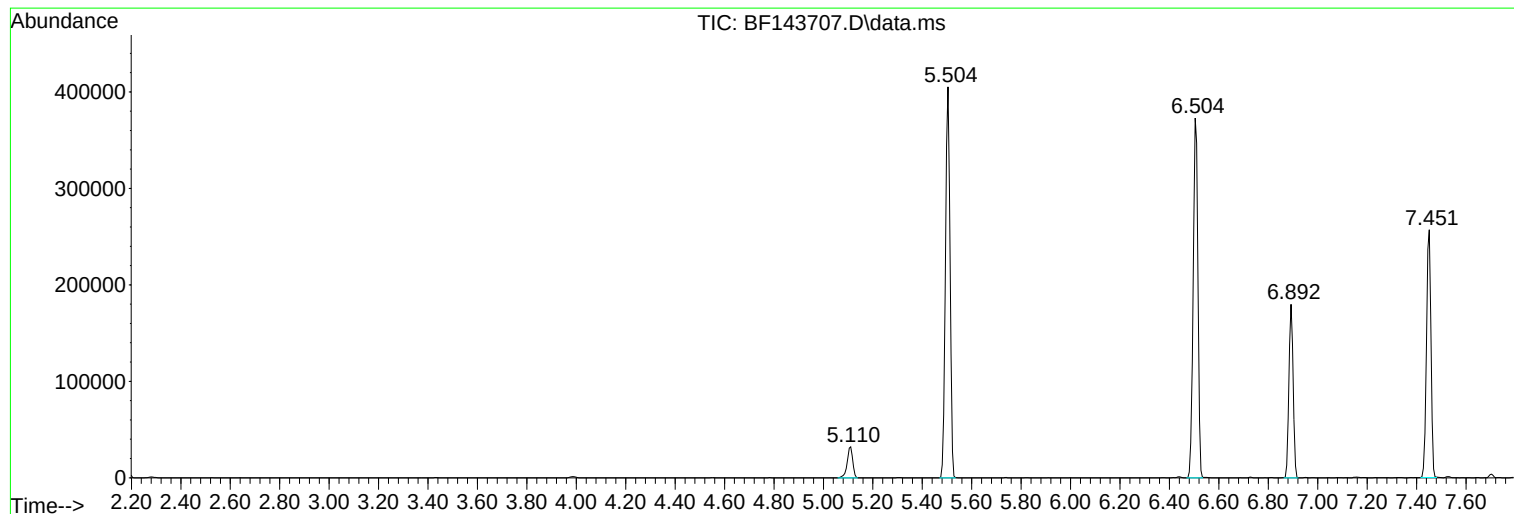
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Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
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ALS Vial : 8 Sample Multiplier: 1

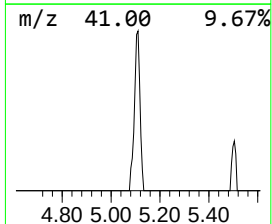
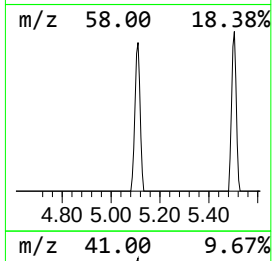
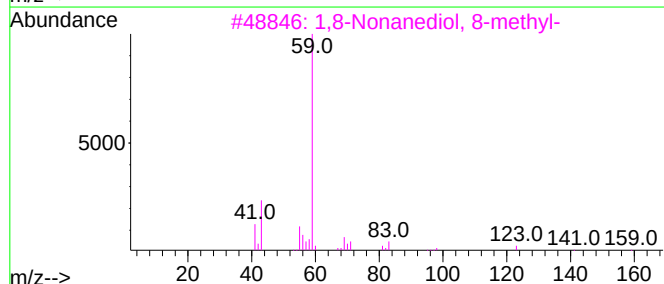
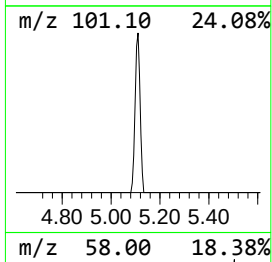
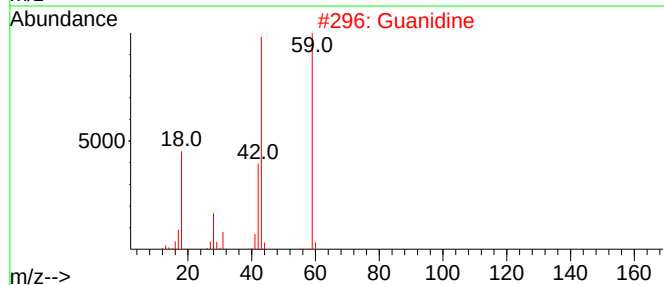
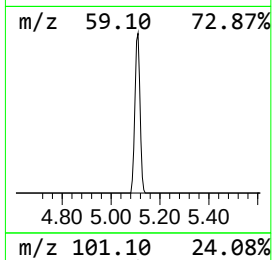
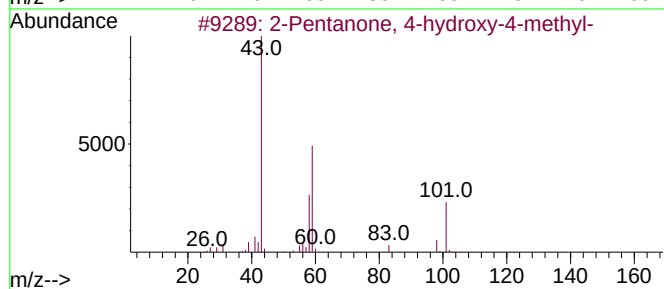
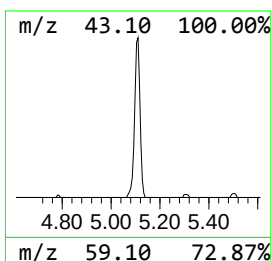
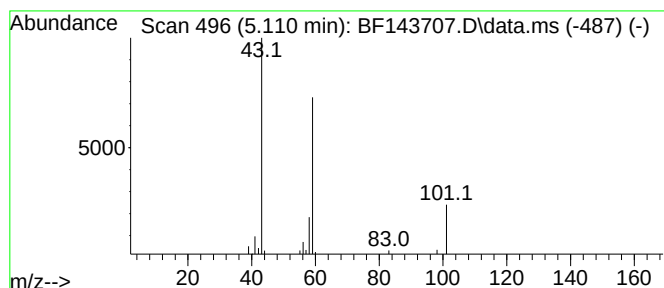
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.110	4.63 ng	49421	1,4-Dichlorobenzene-d4	6.892	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Guanidine	59	CH5N3	000113-00-8	43
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Acetone	58	C3H6O	000067-64-1	9



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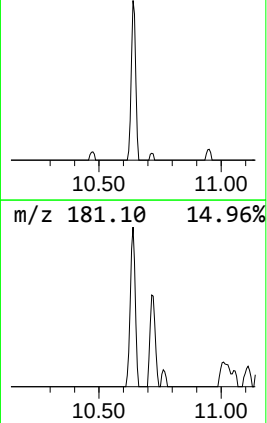
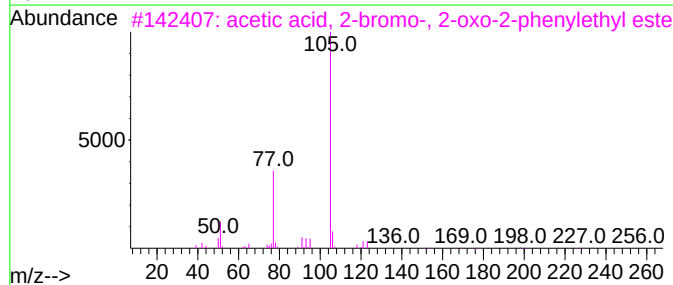
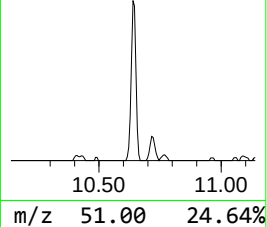
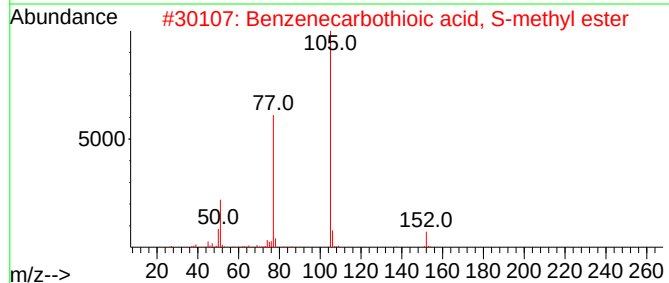
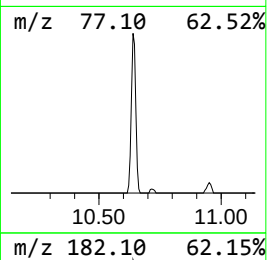
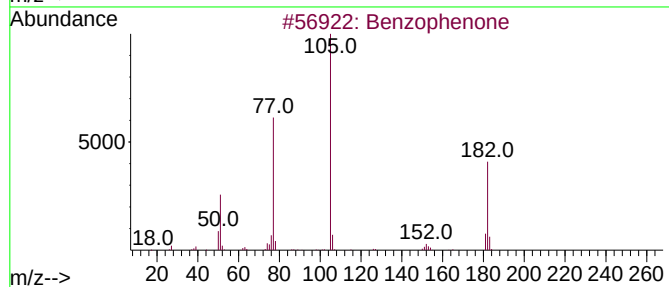
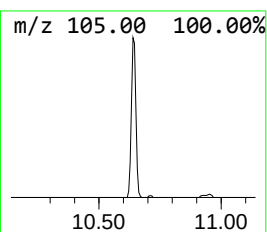
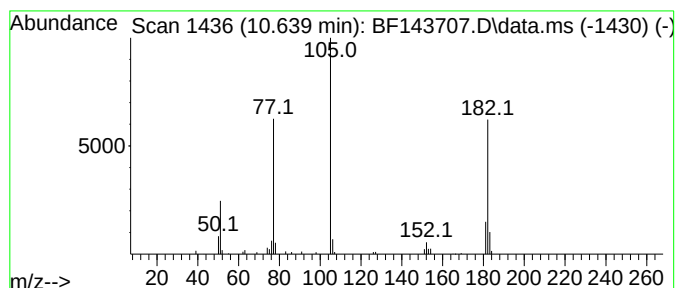
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.639	5.38 ng	84071	Acenaphthene-d10	9.928

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	46
3			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	43
4			2-Benzoyl-4-bromo-5-methyl-1,2-o...	281	C11H8BrNO3	1000444-92-3	43
5			.beta.-Benzilmonoxime	225	C14H11NO2	014090-77-8	43



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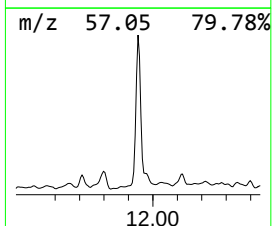
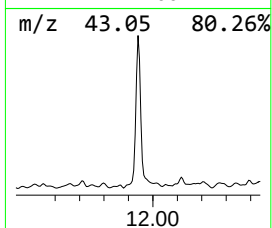
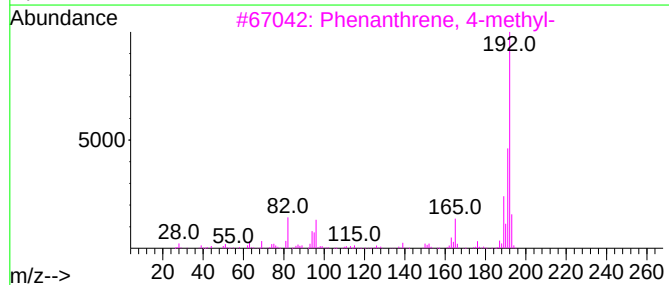
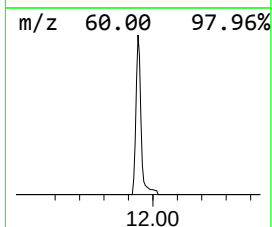
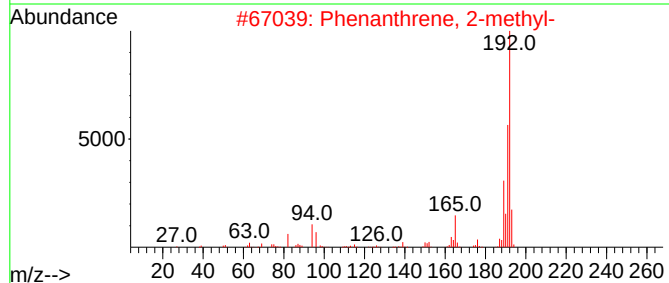
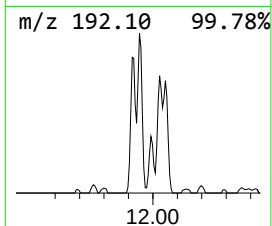
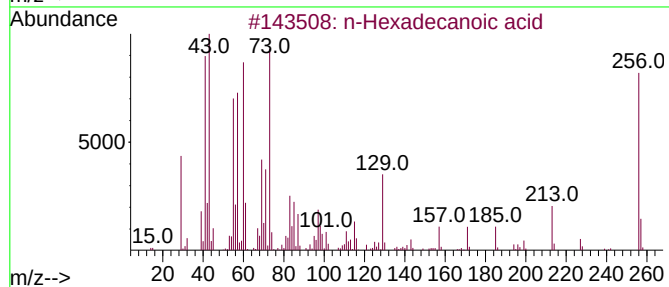
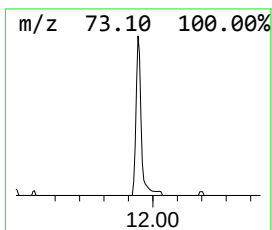
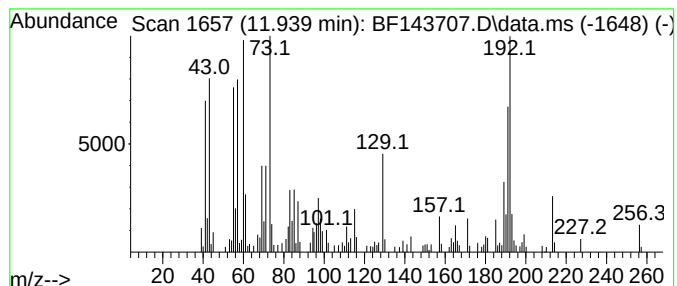
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.939	14.52 ng	191686	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	97
2	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	60
3	Phenanthrene, 4-methyl-		192	C15H12	000832-64-4	60
4	Anthracene, 2-methyl-		192	C15H12	000613-12-7	56
5	Pentadecanoic acid		242	C15H30O2	001002-84-2	55



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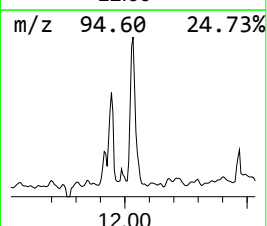
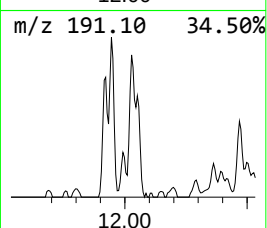
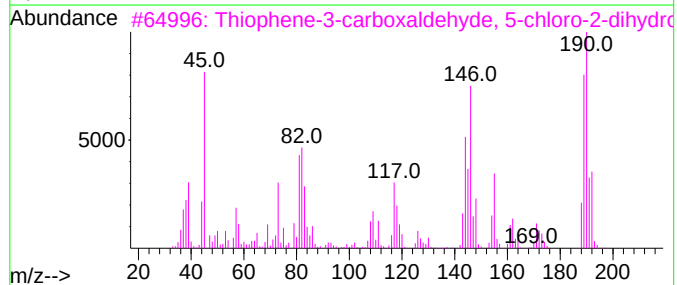
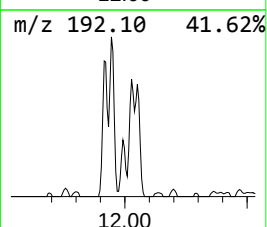
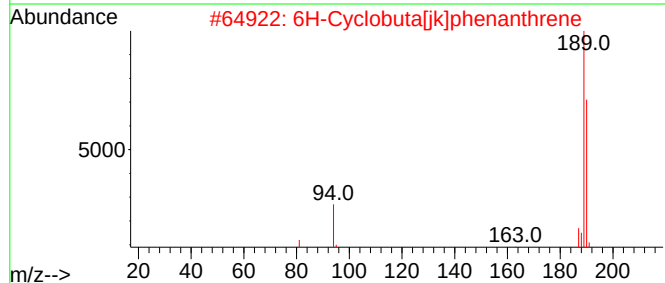
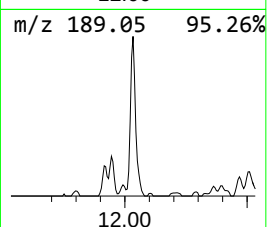
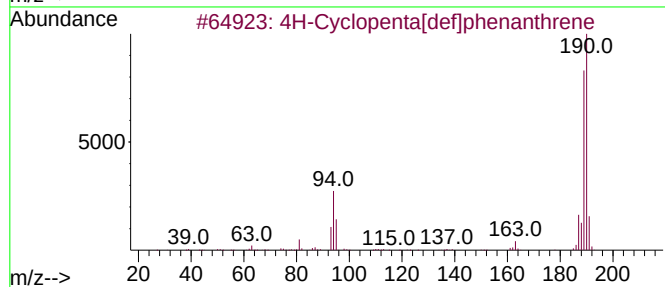
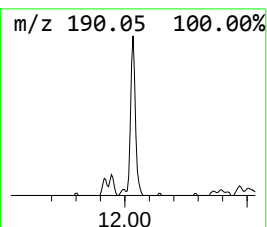
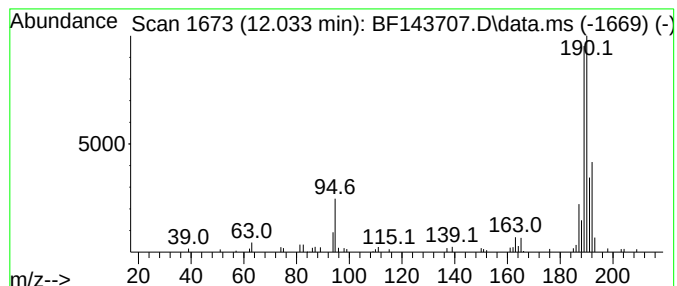
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OR-03-090925

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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 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.033	7.71 ng	101802	Phenanthrene-d10	11.416	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
4	Methyl diselenide	190	C2H6Se2	007101-31-7	43
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
Operator : RC/JU
Sample : Q3055-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

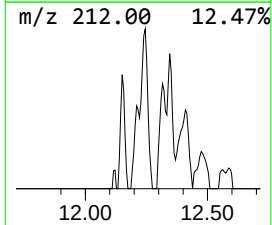
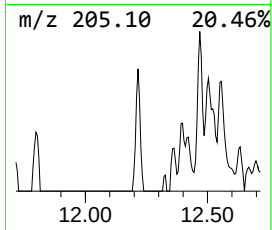
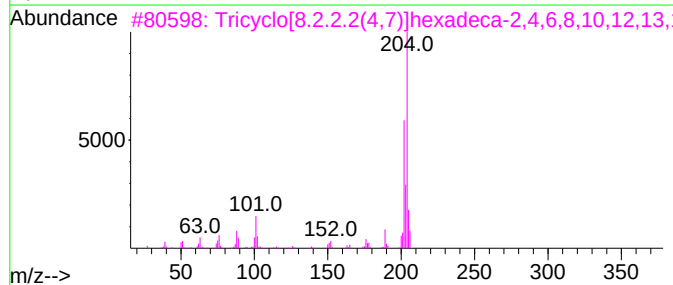
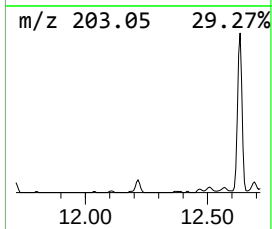
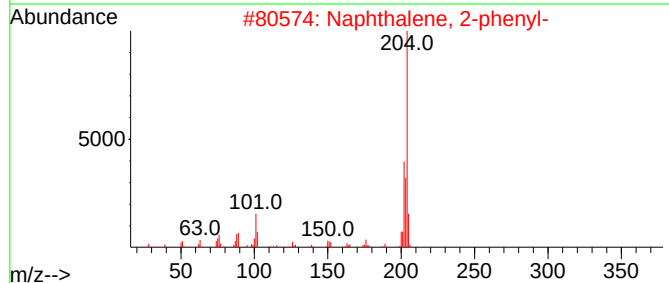
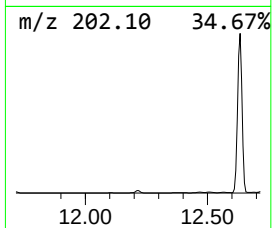
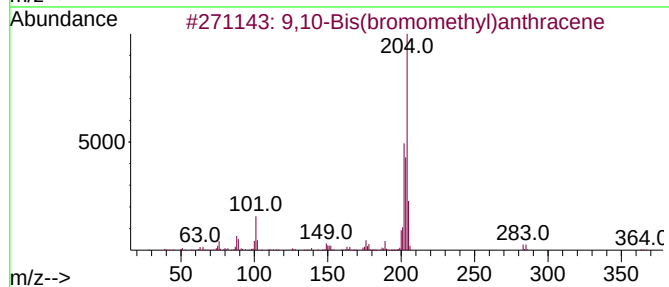
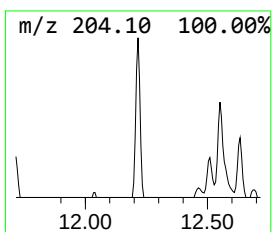
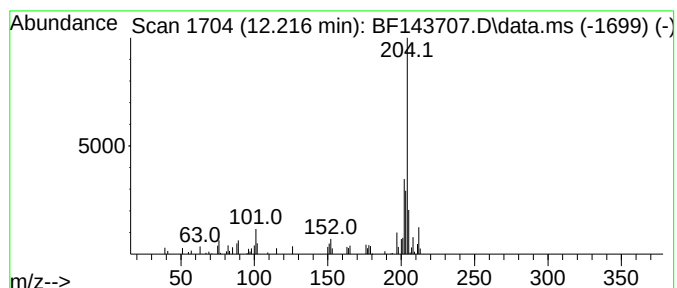
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 9,10-Bis(bromomethyl)anthra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.216	3.93 ng	51918	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9,10-Bis(bromomethyl)anthracene		362	C16H12Br2	034373-96-1	74
2	Naphthalene, 2-phenyl-		204	C16H12	000612-94-2	70
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...		204	C16H12	006572-60-7	70
4	[1,1'-Biphenyl]-4,4'-dicarbonitrile		204	C14H8N2	001591-30-6	49
5	[1,1'-Biphenyl]-2,2'-dicarbonitrile		204	C14H8N2	004341-02-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
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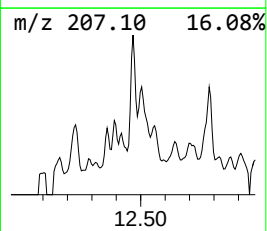
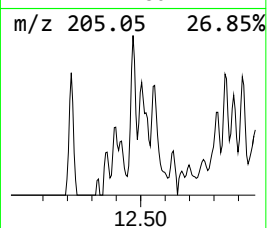
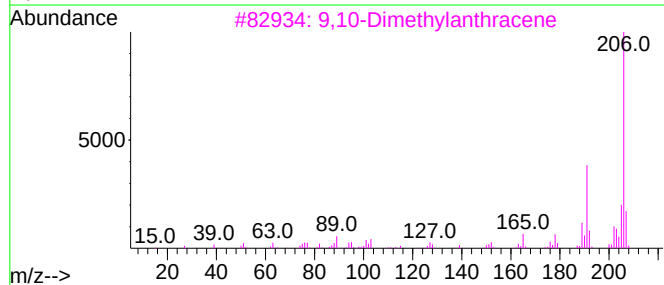
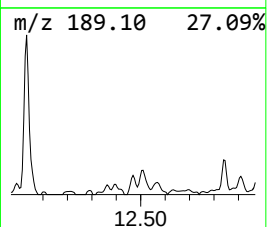
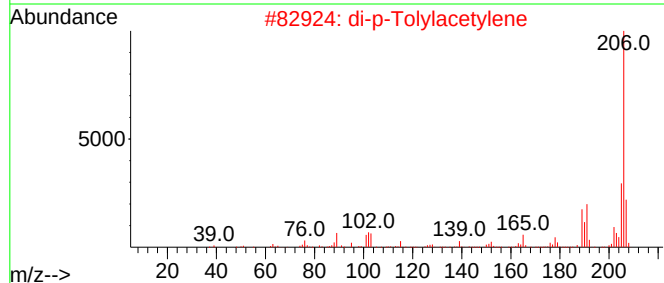
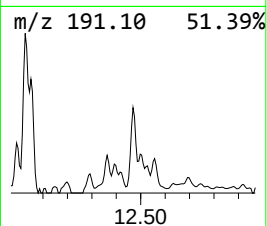
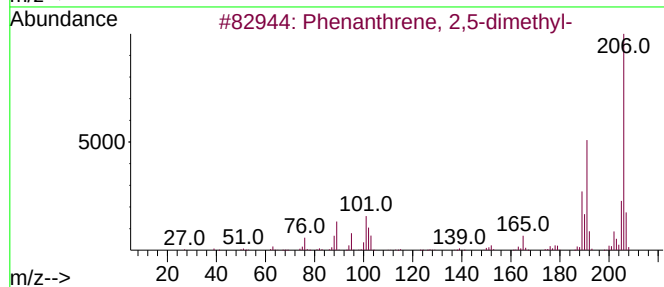
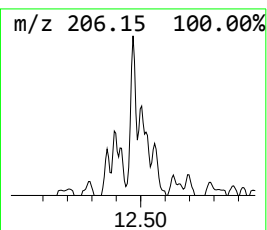
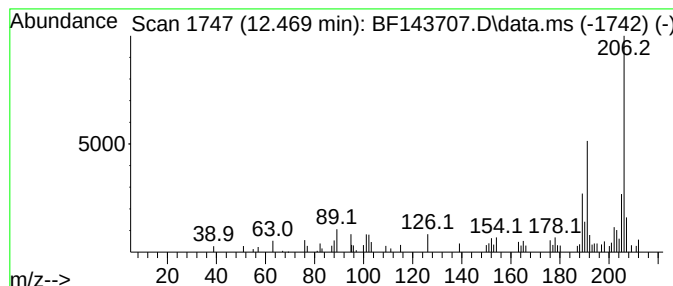
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.469	2.46 ng	32530	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	91
2	di-p-Tolylacetylene		206	C16H14	002789-88-0	91
3	9,10-Dimethylantracene		206	C16H14	000781-43-1	90
4	1H-Indene, 2-methyl-3-phenyl-		206	C16H14	035099-60-6	89
5	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
Operator : RC/JU
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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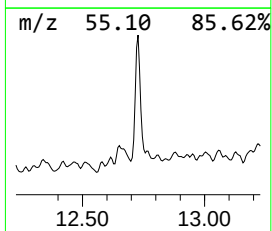
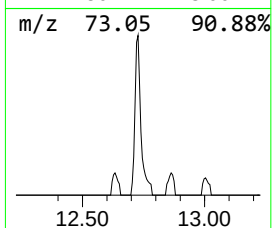
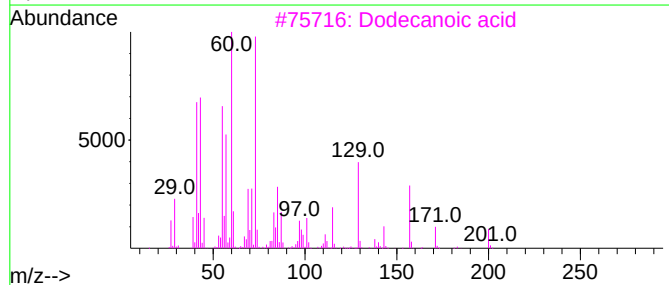
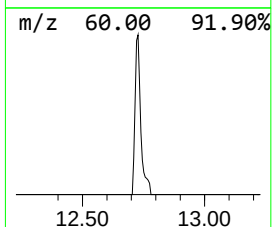
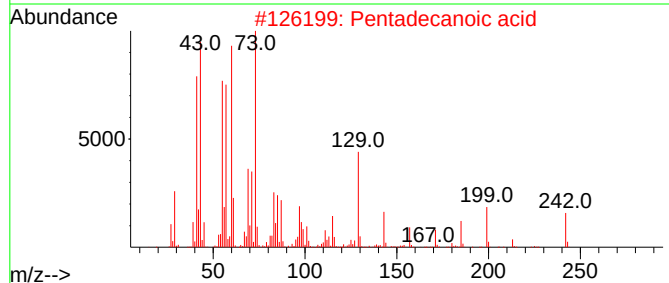
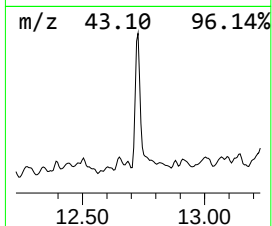
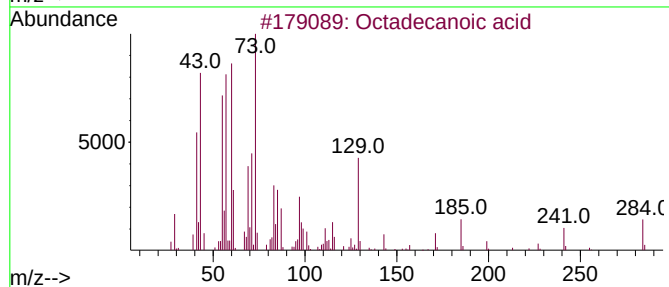
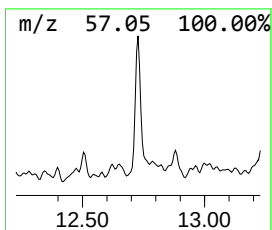
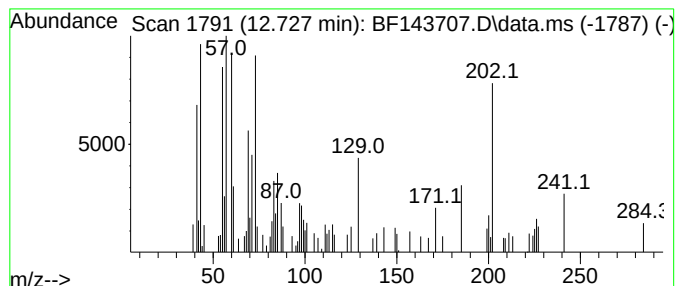
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.727	3.47 ng	45795	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	49
3			Dodecanoic acid	200	C12H24O2	000143-07-7	46
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
Operator : RC/JU
Sample : Q3055-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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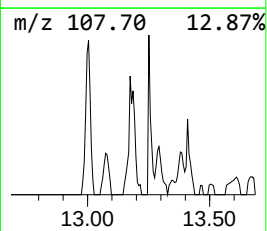
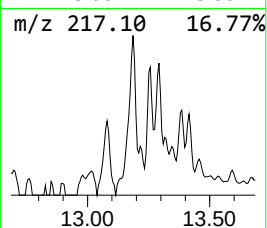
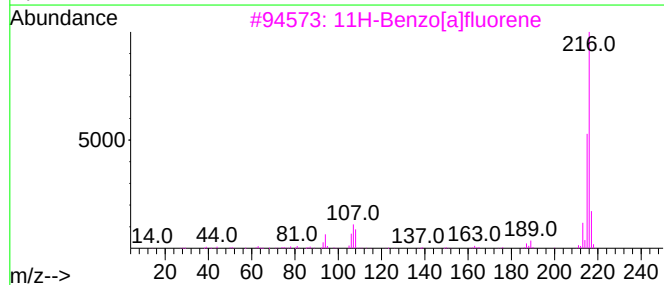
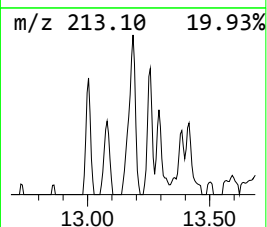
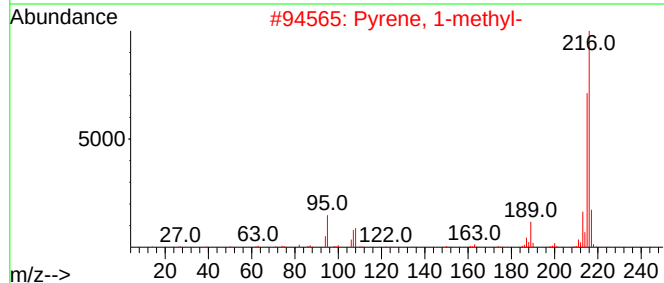
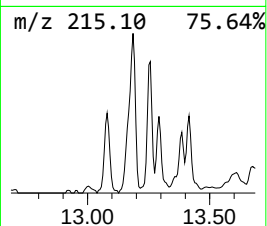
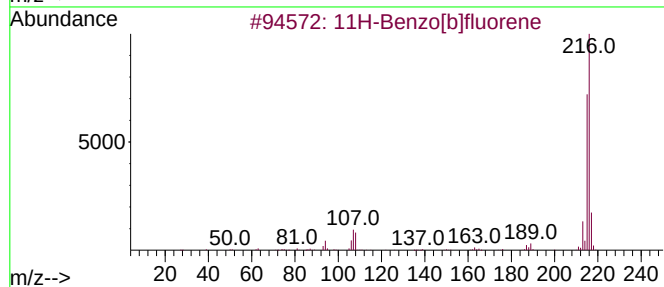
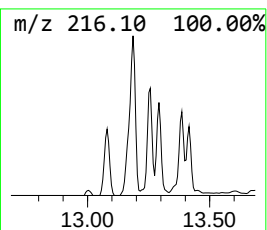
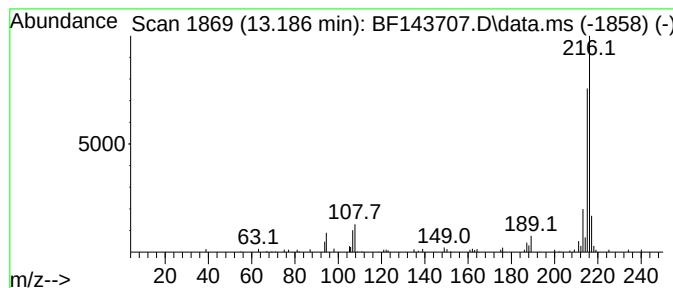
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.186	3.14 ng	80820	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
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Misc :
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Instrument :
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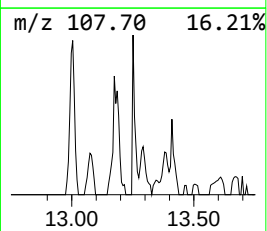
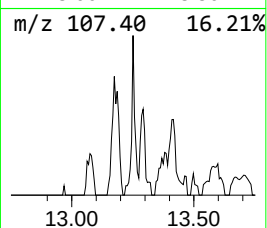
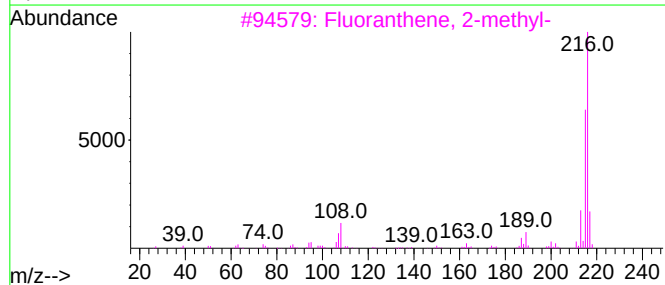
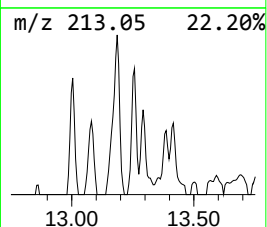
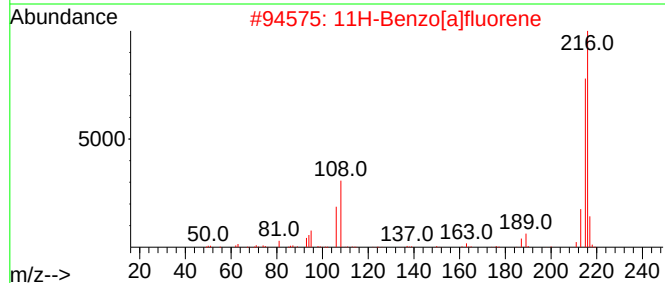
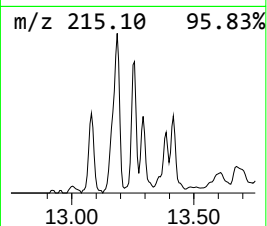
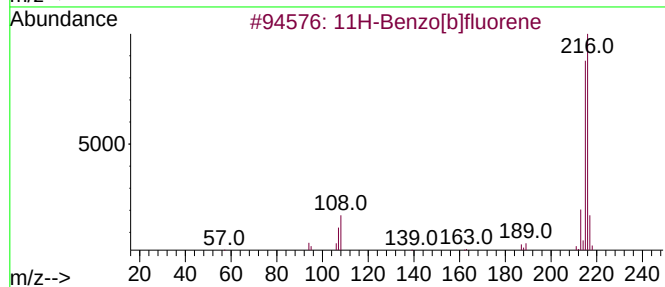
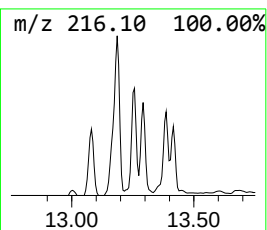
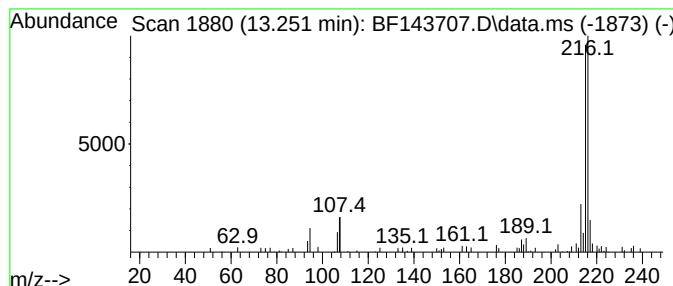
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.251	2.04 ng	52582	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	64
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
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ALS Vial : 8 Sample Multiplier: 1

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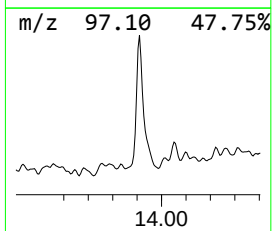
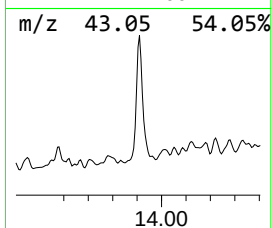
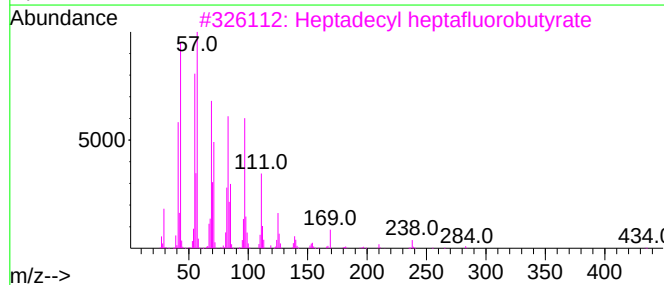
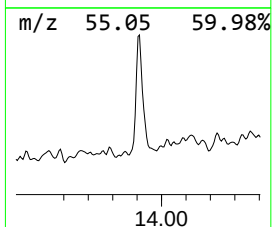
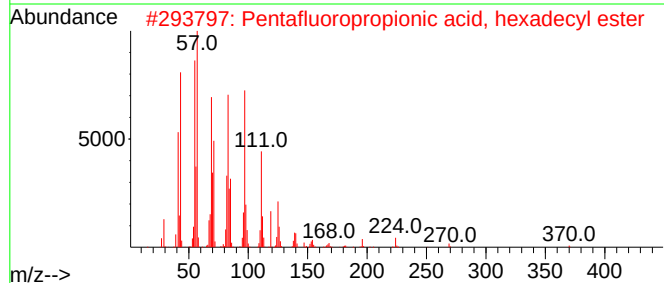
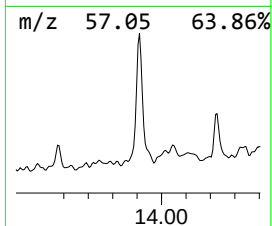
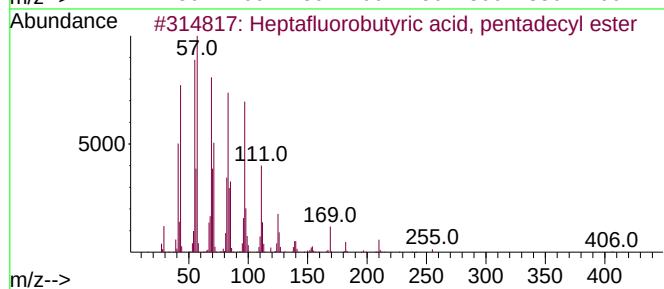
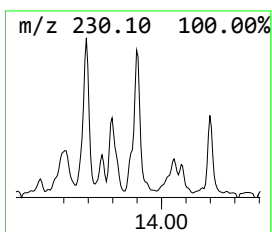
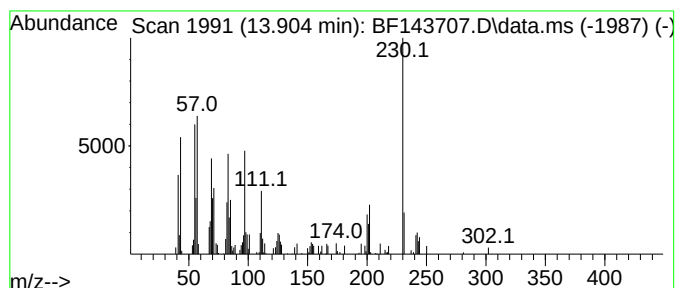
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptafluorobutyric acid, pe... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	2.40 ng	61665	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	80
2			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	46
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	46
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	42
5			3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
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ALS Vial : 8 Sample Multiplier: 1

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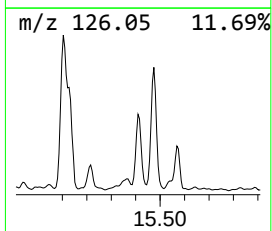
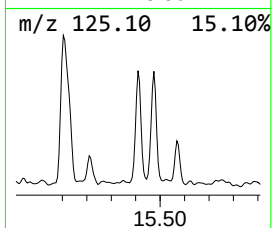
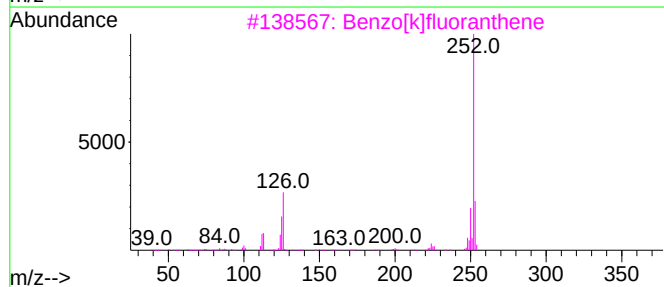
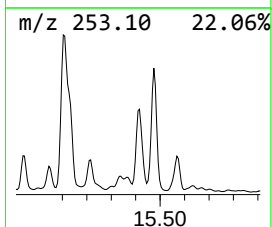
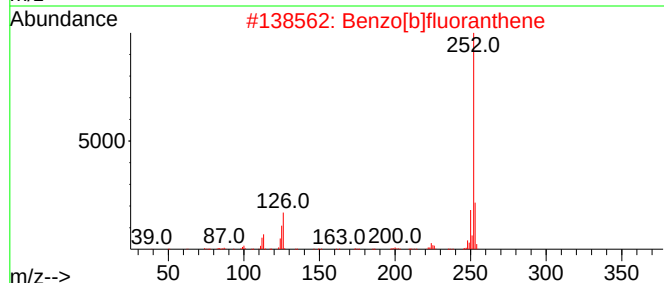
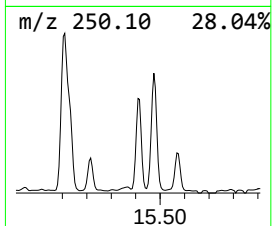
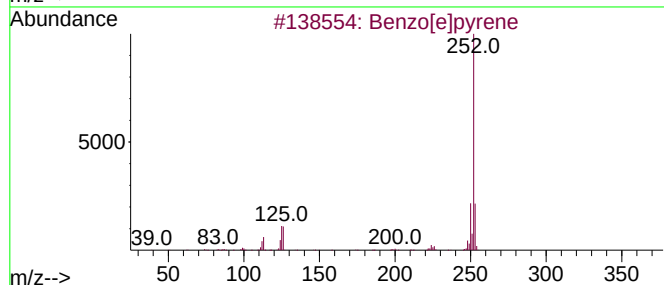
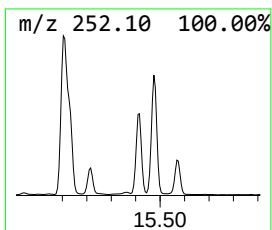
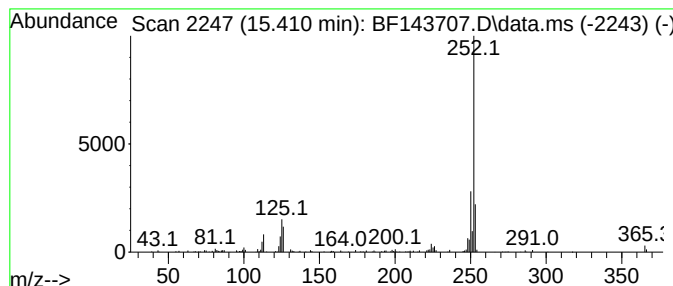
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.410	10.38 ng	107736	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	93
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	91
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091125\
Data File : BF143707.D
Acq On : 11 Sep 2025 15:32
Operator : RC/JU
Sample : Q3055-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-03-090925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.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
2-Pentanone, 4-...	5.110	4.6	ng	49421	1	6.892	213572	20.0
Benzophenone	10.639	5.4	ng	84071	3	9.928	312670	20.0
n-Hexadecanoic ...	11.939	14.5	ng	191686	4	11.416	264082	20.0
4H-Cyclopenta[d...]	12.033	7.7	ng	101802	4	11.416	264082	20.0
9,10-Bis(bromom...	12.216	3.9	ng	51918	4	11.416	264082	20.0
Phenanthrene, 2...	12.469	2.5	ng	32530	4	11.416	264082	20.0
Octadecanoic acid	12.727	3.5	ng	45795	4	11.416	264082	20.0
11H-Benzo[b]flu...	13.186	3.1	ng	80820	5	14.057	514409	20.0
11H-Benzo[a]flu...	13.251	2.0	ng	52582	5	14.057	514409	20.0
Heptafluorobuty...	13.904	2.4	ng	61665	5	14.057	514409	20.0
Benzo[e]pyrene	15.410	10.4	ng	107736	6	15.539	207623	20.0
unknown16.268	16.268	3.1	ng	32080	6	15.539	207623	20.0