

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
 Data File : BF139568.D
 Acq On : 27 Sep 2024 01:42
 Operator : RC/JU
 Sample : P4197-01 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-092524

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: BF139568.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.134	3	7	22	rVB	430717	639234	66.70%	5.158%
2	3.381	214	219	230	rBV	3305	11216	1.17%	0.091%
3	3.934	304	313	320	rVB	6493	11450	1.19%	0.092%
4	5.075	498	507	512	rVB	18852	25918	2.70%	0.209%
5	5.487	566	577	593	rVB	281743	375075	39.14%	3.026%
6	6.498	744	749	757	rBV	305096	387706	40.46%	3.128%
7	6.704	780	784	789	rBV3	8043	11483	1.20%	0.093%
8	6.875	808	813	818	rBV	435677	525215	54.81%	4.238%
9	7.433	899	908	913	rBV	200689	260415	27.17%	2.101%
10	7.516	917	922	926	rVB5	6146	10119	1.06%	0.082%
11	7.681	945	950	957	rVB5	11706	22471	2.34%	0.181%
12	7.775	957	966	967	rBV5	4881	9737	1.02%	0.079%
13	7.892	979	986	989	rBV4	10287	16685	1.74%	0.135%
14	8.157	1017	1031	1038	rBV	553089	748591	78.11%	6.040%
15	8.445	1075	1080	1082	rBV4	11545	18300	1.91%	0.148%
16	8.539	1093	1096	1099	rBV2	9053	9787	1.02%	0.079%
17	8.657	1113	1116	1120	rBV	24034	30720	3.21%	0.248%
18	8.745	1127	1131	1134	rBV2	18861	24725	2.58%	0.200%
19	8.816	1140	1143	1149	rVB4	11219	16698	1.74%	0.135%
20	8.869	1149	1152	1155	rVB	22975	24973	2.61%	0.202%
21	8.975	1164	1170	1176	rBV3	39225	69529	7.26%	0.561%
22	9.104	1188	1192	1197	rBV5	10523	18263	1.91%	0.147%
23	9.175	1201	1204	1209	rVV4	10749	15491	1.62%	0.125%
24	9.227	1209	1213	1218	rBV	357578	428661	44.73%	3.459%
25	9.310	1223	1227	1230	rBV	35399	47846	4.99%	0.386%
26	9.380	1235	1239	1245	rVB	15497	22661	2.36%	0.183%
27	9.492	1253	1258	1263	rBV2	19662	33373	3.48%	0.269%
28	9.569	1264	1271	1273	rVV2	33496	62635	6.54%	0.505%
29	9.627	1277	1281	1284	rVV3	32358	53936	5.63%	0.435%
30	9.686	1288	1291	1296	rVB2	23117	31178	3.25%	0.252%
31	9.769	1301	1305	1309	rBV	60586	81998	8.56%	0.662%
32	9.839	1313	1317	1321	rVB	49329	62942	6.57%	0.508%
33	9.910	1324	1329	1333	rBV	557876	689747	71.97%	5.565%
34	10.010	1343	1346	1351	rVB3	16048	22730	2.37%	0.183%
35	10.069	1352	1356	1358	rBV4	12702	20890	2.18%	0.169%
36	10.116	1361	1364	1367	rVB2	28888	37809	3.95%	0.305%

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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	10.157	1367	1371	1375	rBV4	18402	23454	2.45%	0.189%
38	10.227	1379	1383	1385	rBV3	15480	21794	2.27%	0.176%
39	10.333	1395	1401	1405	rBV	75556	113361	11.83%	0.915%
40	10.457	1418	1422	1429	rBV	53427	73511	7.67%	0.593%
41	10.557	1434	1439	1444	rVB	34197	58462	6.10%	0.472%
42	10.610	1445	1448	1451	rBV3	13993	16665	1.74%	0.134%
43	10.698	1457	1463	1467	rBV	155543	208612	21.77%	1.683%
44	10.810	1478	1482	1491	rVB	100349	171635	17.91%	1.385%
45	11.004	1511	1515	1518	rBV2	22241	37164	3.88%	0.300%
46	11.133	1533	1537	1539	rBV2	14149	21913	2.29%	0.177%
47	11.257	1554	1558	1561	rBV	77987	102660	10.71%	0.828%
48	11.286	1561	1563	1569	rVB2	53574	72409	7.56%	0.584%
49	11.398	1577	1582	1585	rBV	503978	748832	78.14%	6.042%
50	11.469	1591	1594	1598	rVB	43424	49949	5.21%	0.403%
51	11.504	1598	1600	1604	rVB	13800	14097	1.47%	0.114%
52	11.633	1618	1622	1626	rBV2	24284	34289	3.58%	0.277%
53	11.686	1627	1631	1635	rBV2	65130	82273	8.59%	0.664%
54	11.786	1644	1648	1655	rVB	79885	103656	10.82%	0.836%
55	11.921	1663	1671	1676	rBV2	74008	155757	16.25%	1.257%
56	11.974	1677	1680	1682	rVV2	20359	26933	2.81%	0.217%
57	12.010	1682	1686	1693	rVB	75056	133991	13.98%	1.081%
58	12.092	1696	1700	1704	rVV	51546	62739	6.55%	0.506%
59	12.192	1714	1717	1720	rBV	19444	26216	2.74%	0.212%
60	12.374	1745	1748	1750	rBV	11708	13655	1.42%	0.110%
61	12.492	1756	1768	1779	rBV4	269539	664961	69.39%	5.365%
62	12.574	1779	1782	1784	rVV2	33896	39309	4.10%	0.317%
63	12.610	1784	1788	1800	rVV	358768	700870	73.13%	5.655%
64	12.704	1801	1804	1811	rVB	38921	57469	6.00%	0.464%
65	12.839	1820	1827	1833	rBV	298304	441548	46.07%	3.563%
66	12.980	1845	1851	1858	rVB	262085	390720	40.77%	3.153%
67	13.057	1859	1864	1867	rBV4	34200	61396	6.41%	0.495%
68	13.163	1875	1882	1886	rVB2	61721	110641	11.55%	0.893%
69	13.210	1886	1890	1891	rBV	27666	34100	3.56%	0.275%
70	13.233	1891	1894	1897	rVV2	46834	63994	6.68%	0.516%
71	13.268	1897	1900	1907	rVB2	42013	58710	6.13%	0.474%
72	13.363	1913	1916	1918	rBV	17060	18983	1.98%	0.153%
73	13.392	1919	1921	1924	rVB	16548	15973	1.67%	0.129%
74	13.786	1985	1988	1990	rBV	21905	25058	2.61%	0.202%
75	14.033	2024	2030	2041	rBV2	480212	958326	100.00%	7.733%
76	15.080	2203	2208	2216	rBV	182425	399844	41.72%	3.226%

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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-BF091324.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.380	2256	2259	2265	rVB	97979	153759	16.04%	1.241%
78	15.445	2266	2270	2276	rVV	142548	217801	22.73%	1.757%
79	15.509	2276	2281	2294	rVB2	248104	467499	48.78%	3.772%
80	16.980	2526	2531	2539	rVB2	67815	148569	15.50%	1.199%
81	17.427	2603	2607	2620	rVB	50587	109555	11.43%	0.884%

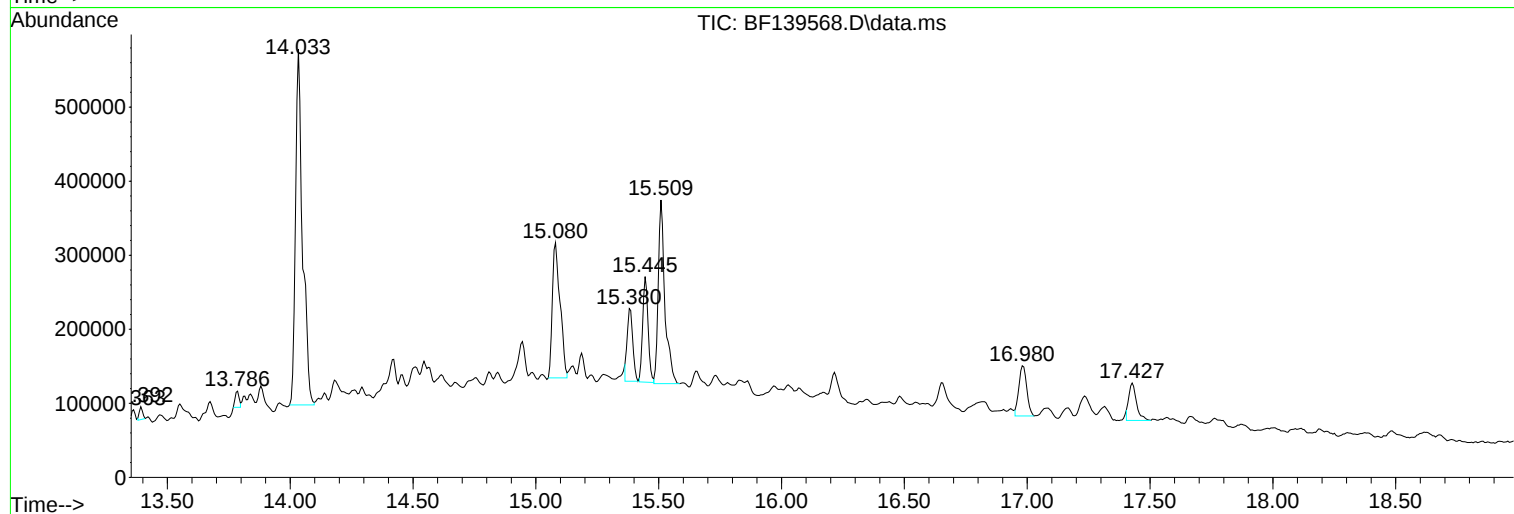
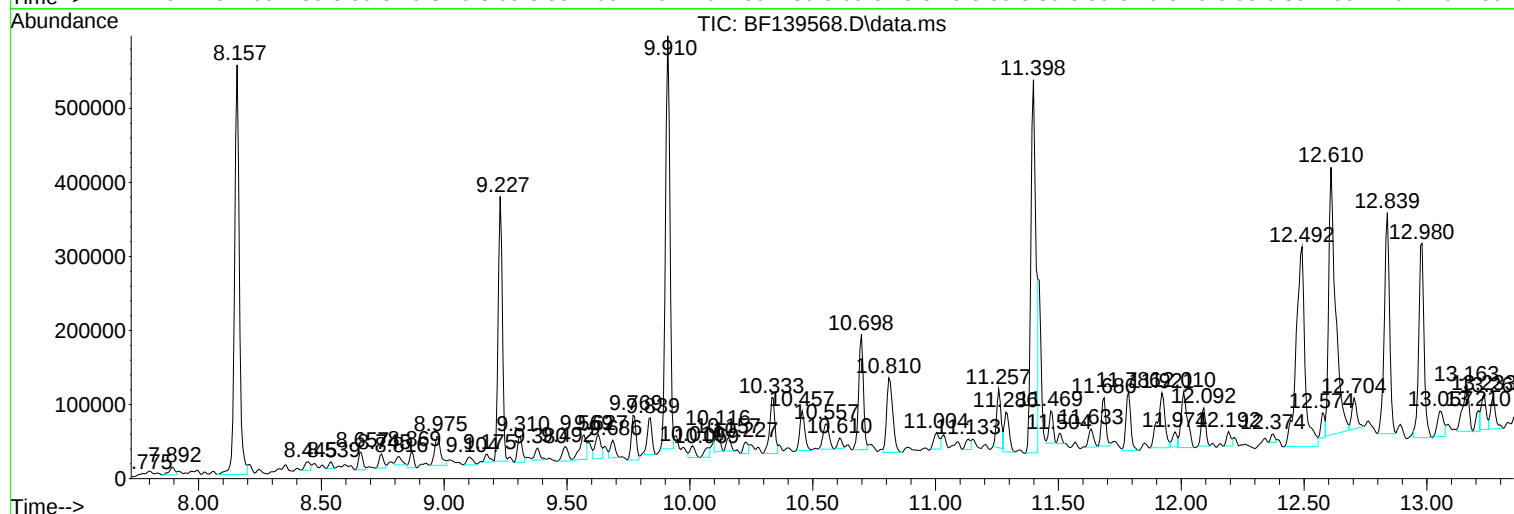
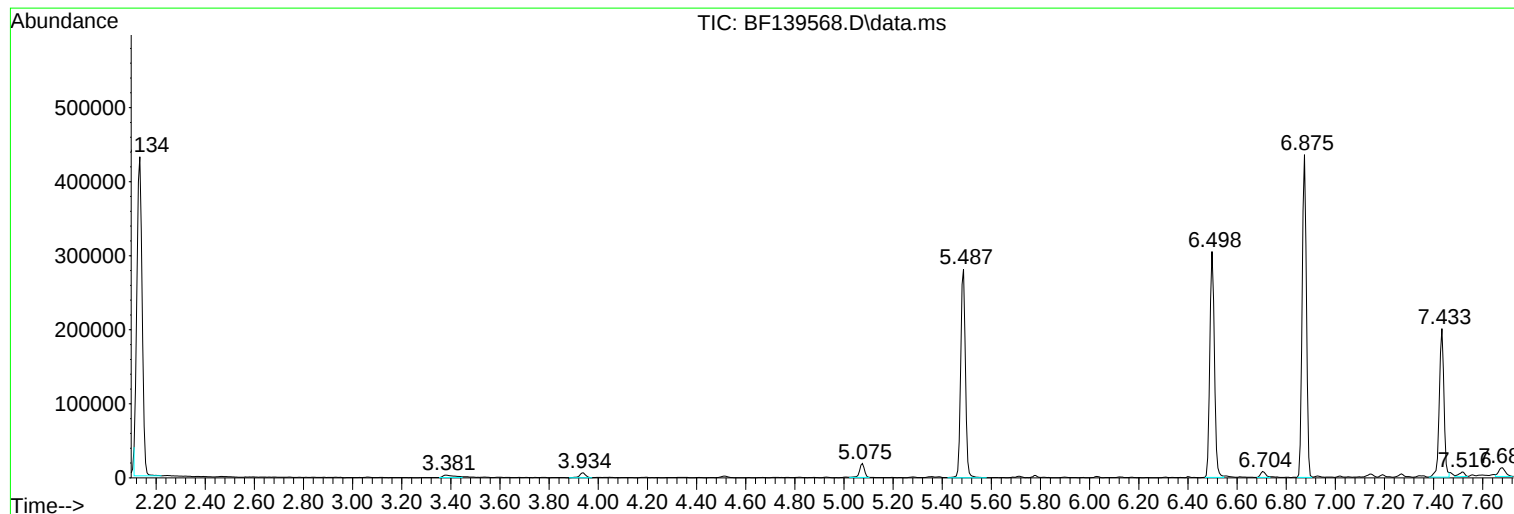
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
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Sample : P4197-01 5X
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Instrument :
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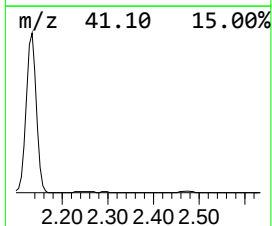
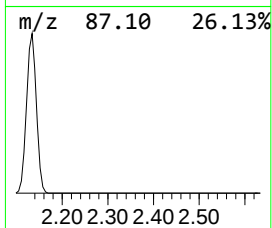
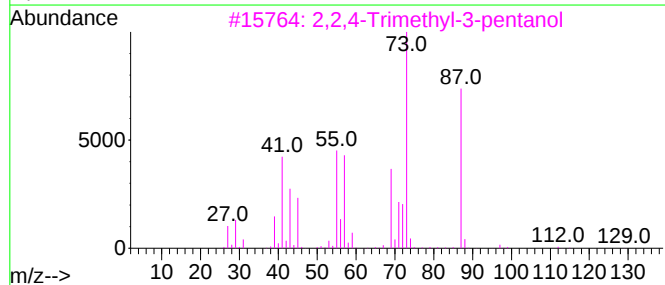
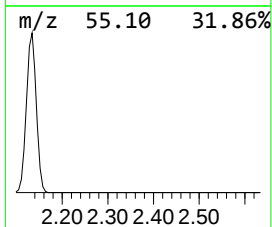
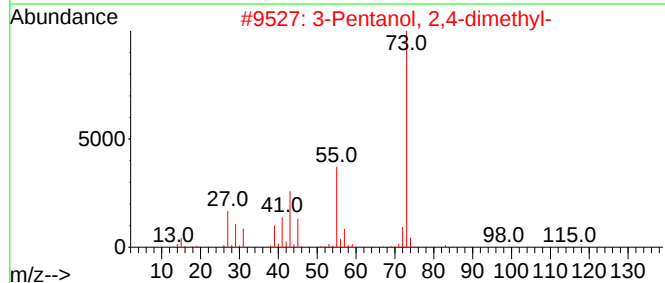
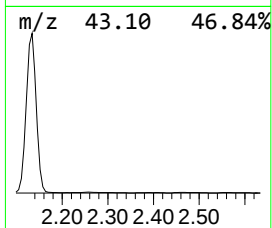
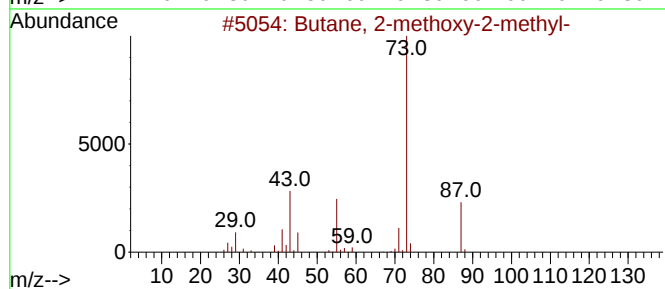
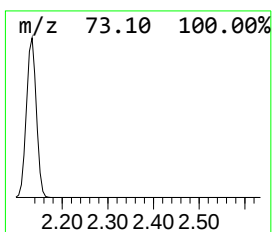
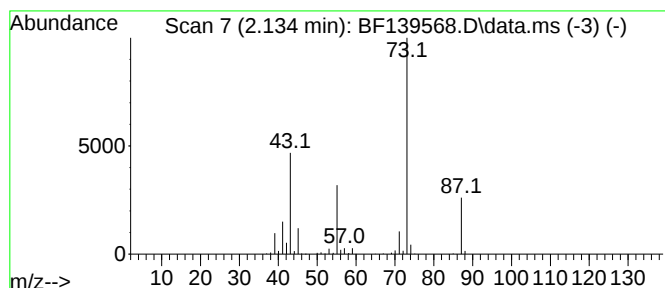
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.134	24.34 ng	639234	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			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	38



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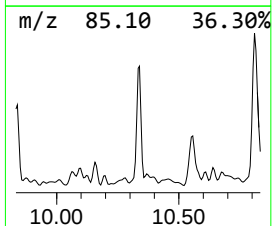
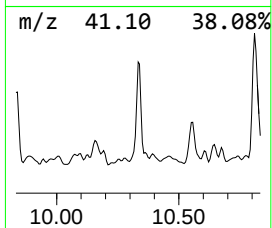
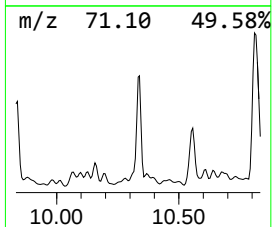
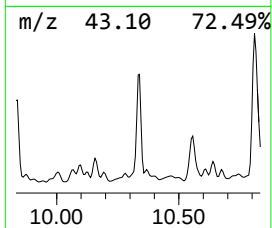
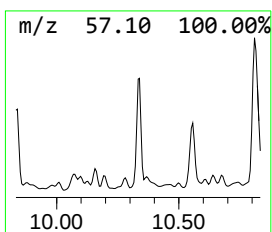
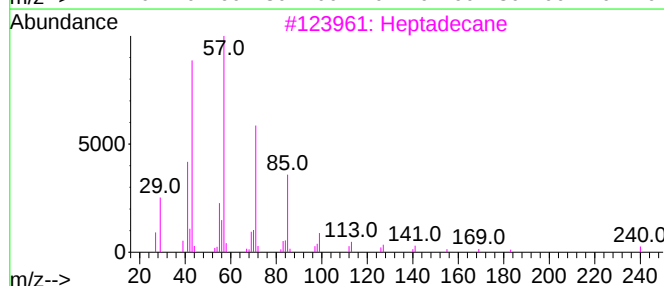
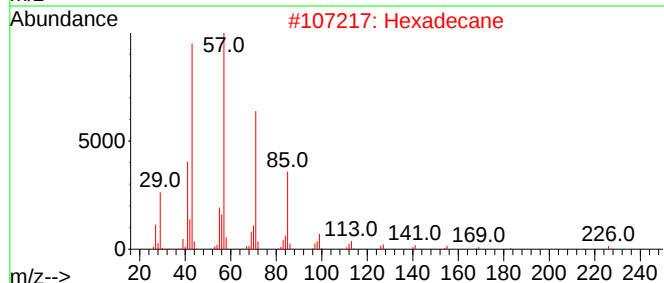
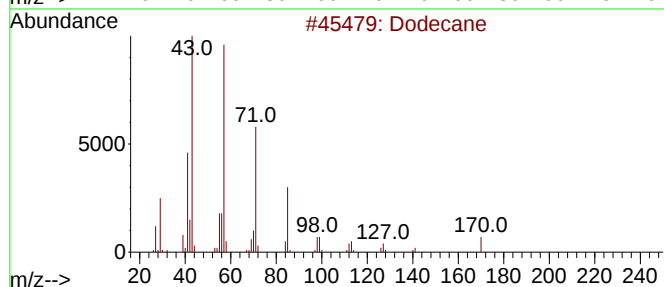
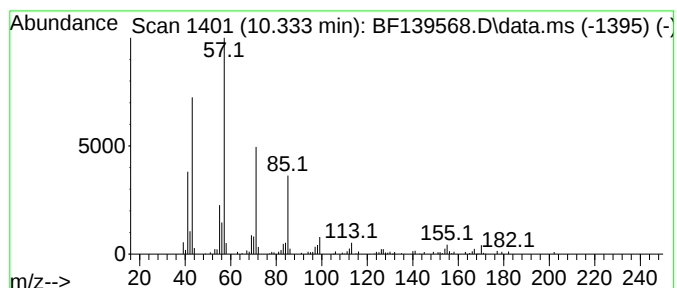
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dodecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.333	3.29 ng	113361	Acenaphthene-d10	9.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	93
2			Hexadecane	226	C16H34	000544-76-3	87
3			Heptadecane	240	C17H36	000629-78-7	86
4			Pentadecane	212	C15H32	000629-62-9	86
5			Tetracosane	338	C24H50	000646-31-1	80



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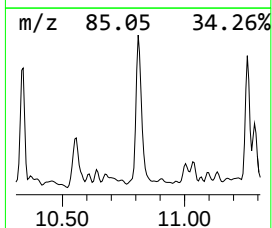
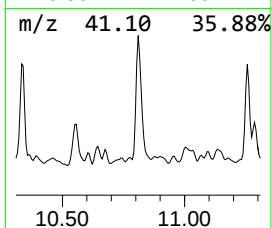
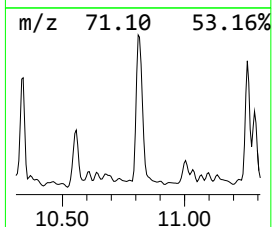
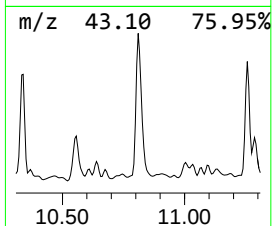
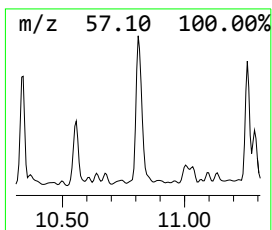
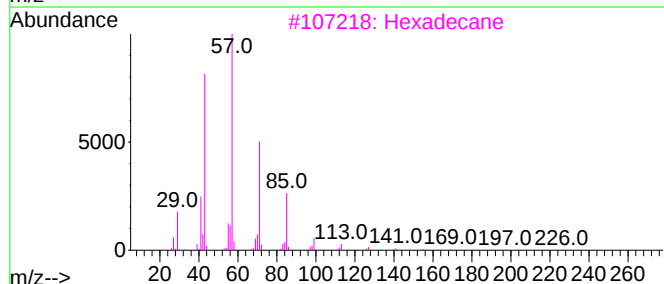
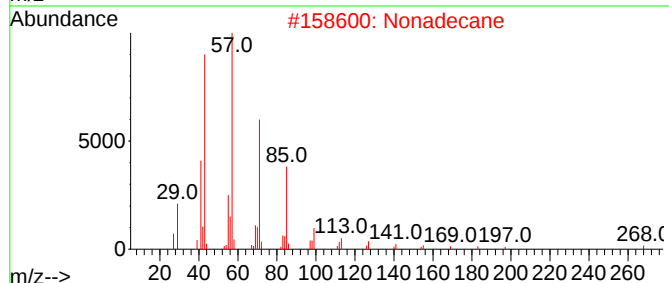
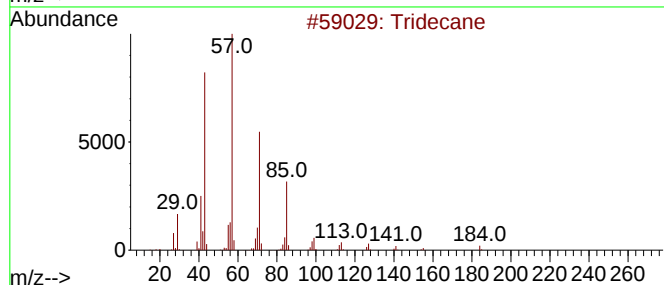
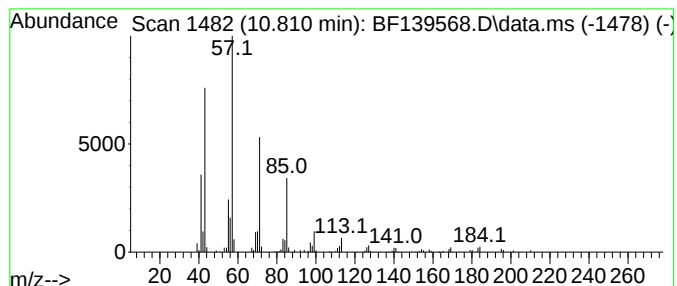
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.810	4.58 ng	171635	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Nonadecane	268	C19H40	000629-92-5	91
3			Hexadecane	226	C16H34	000544-76-3	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Eicosane	282	C20H42	000112-95-8	87



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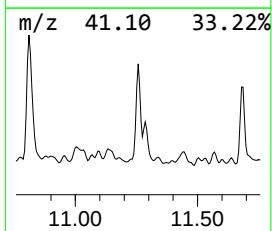
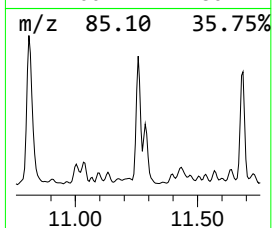
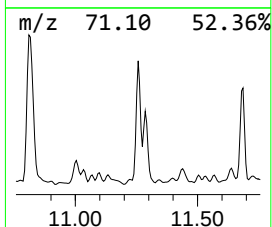
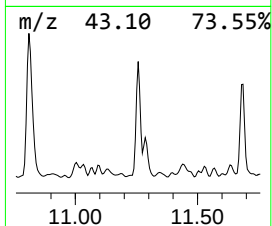
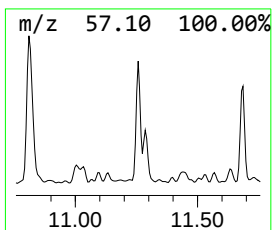
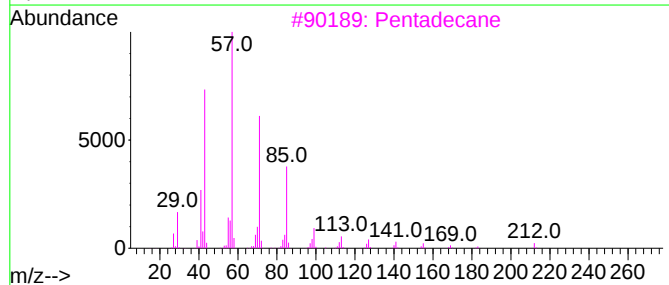
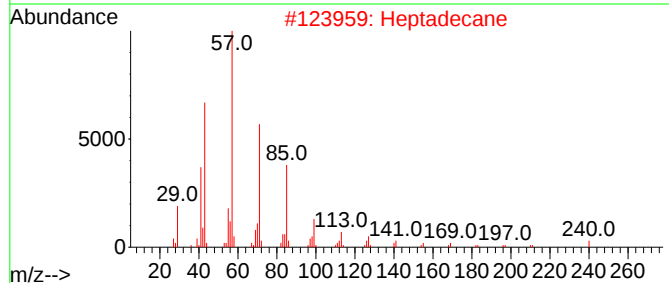
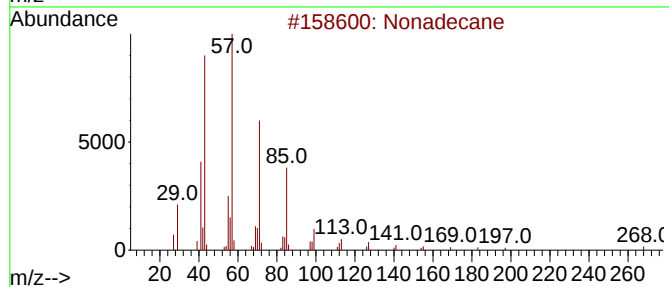
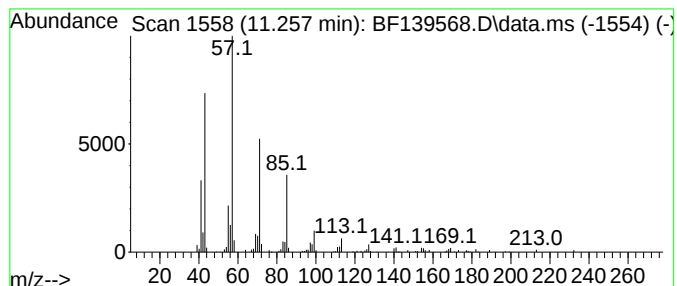
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.257	2.74 ng	102660	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Heptadecane	240	C17H36	000629-78-7	91
3			Pentadecane	212	C15H32	000629-62-9	91
4			Hexadecane	226	C16H34	000544-76-3	90
5			Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-092524

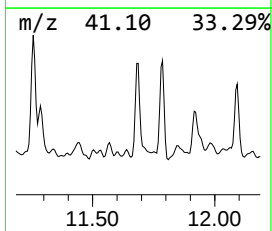
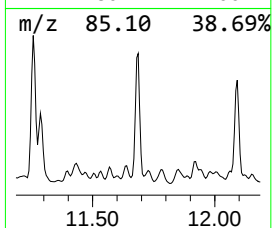
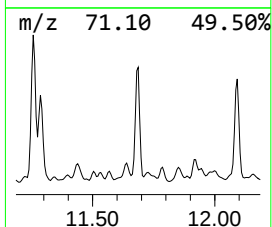
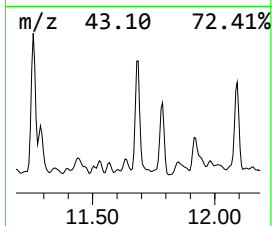
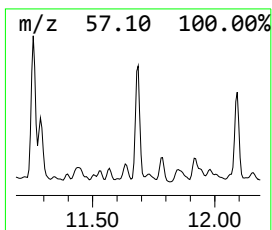
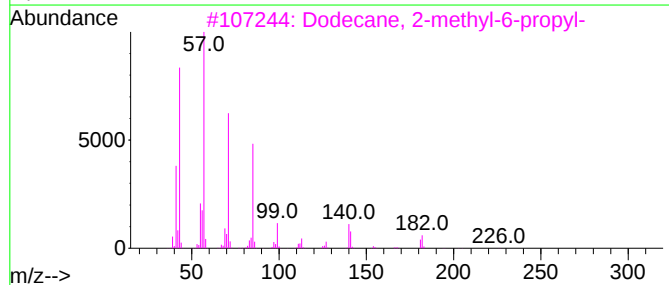
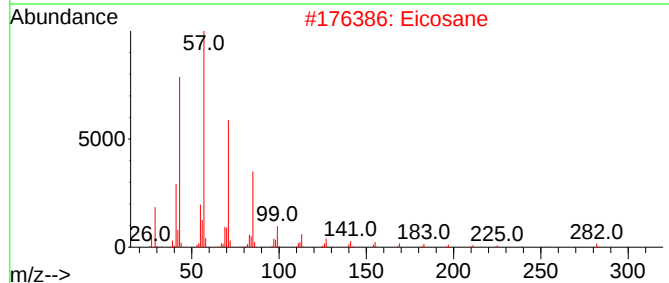
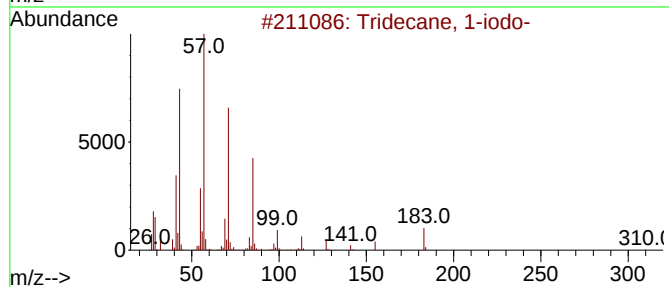
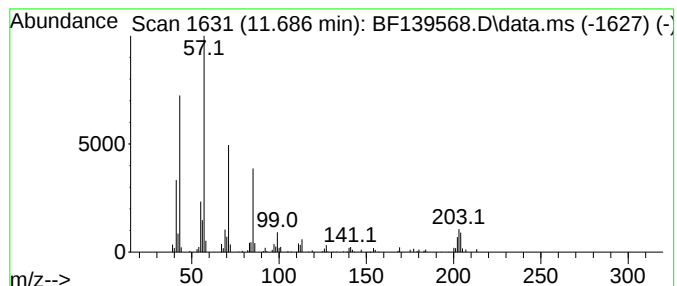
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.686	2.20 ng	82273	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	74
2			Eicosane	282	C20H42	000112-95-8	74
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	72
4			Octacosane	394	C28H58	000630-02-4	72
5			Tridecane, 6-propyl-	226	C16H34	055045-10-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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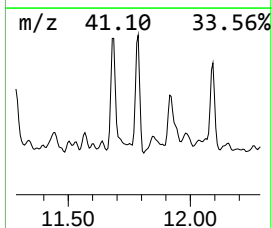
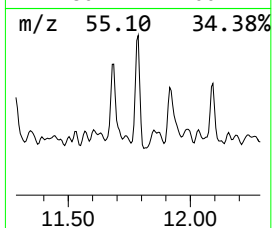
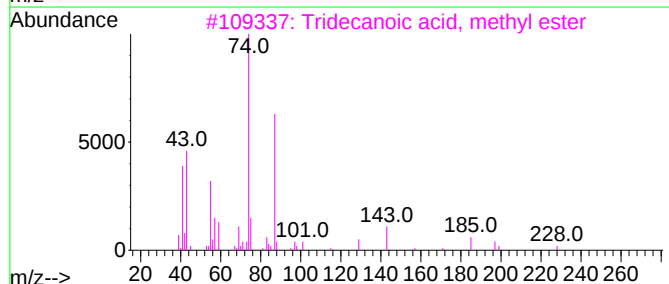
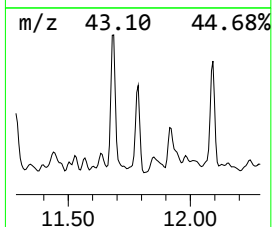
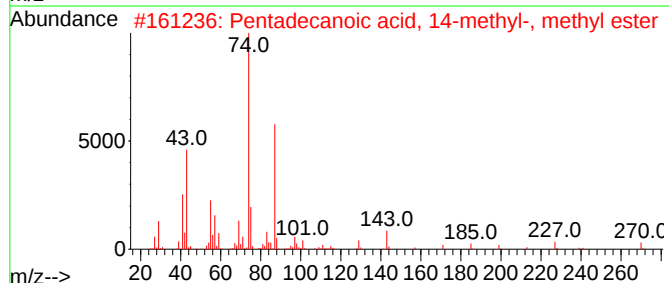
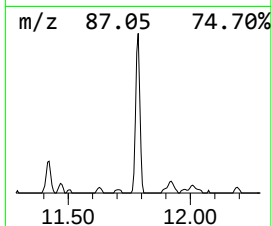
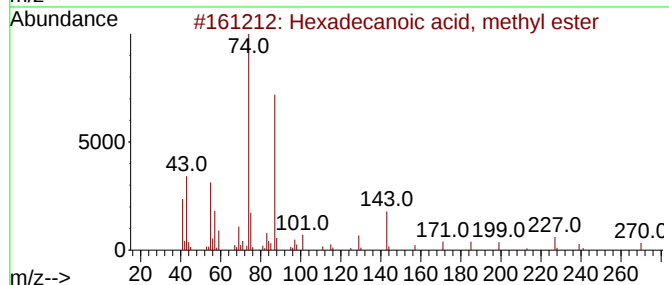
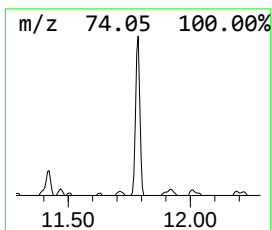
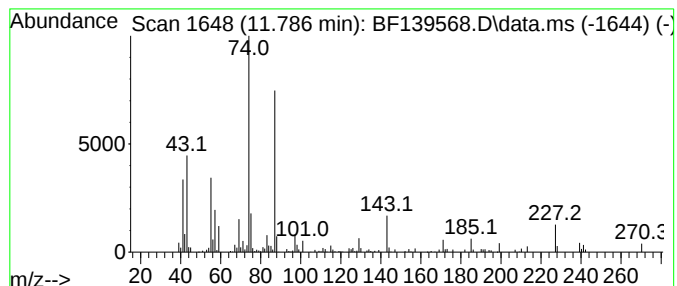
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, methyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.786	2.77 ng	103656	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	95
4			Methyl tetradecanoate	242	C15H30O2	000124-10-7	83
5			Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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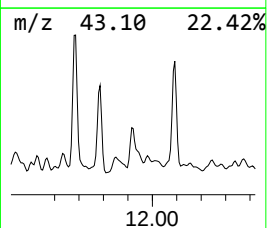
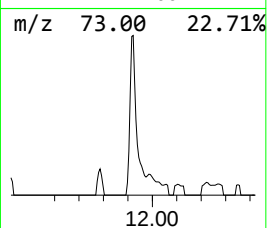
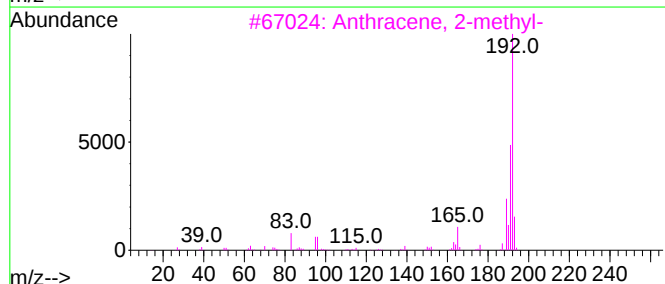
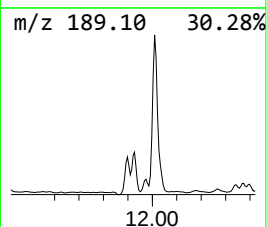
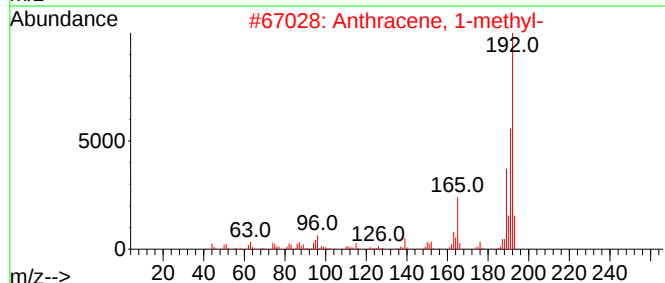
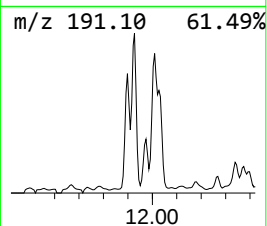
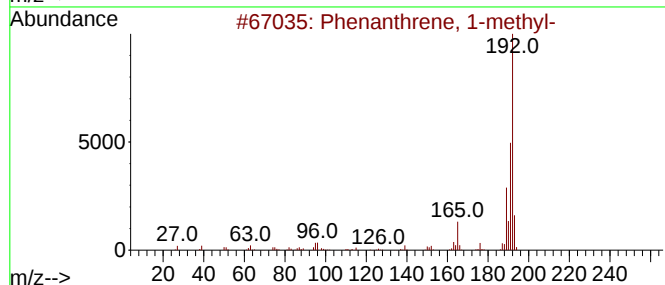
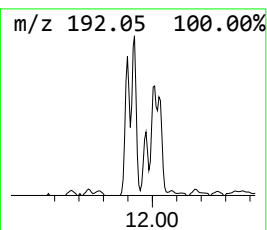
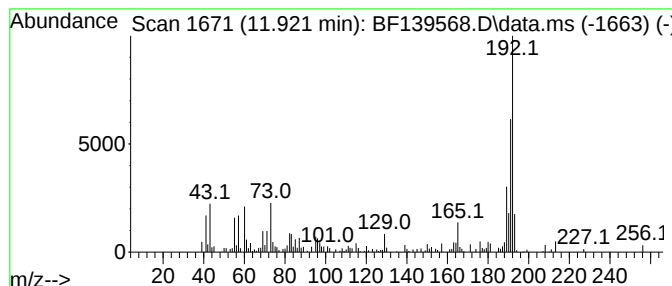
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.921	4.16 ng	155757	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	86
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

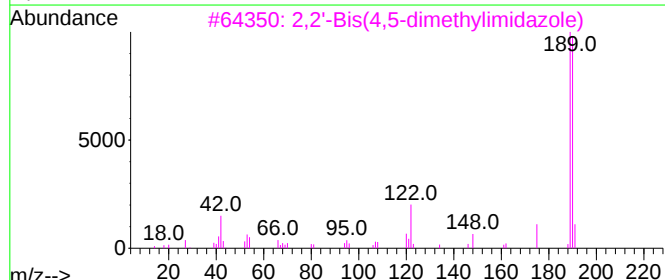
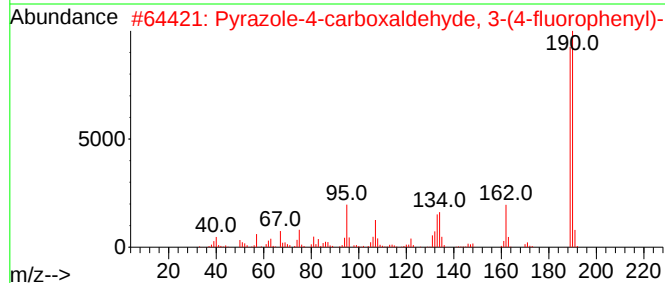
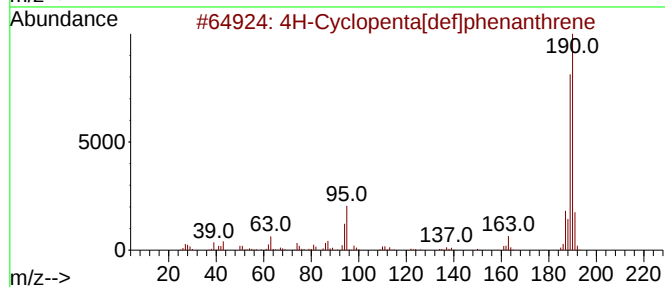
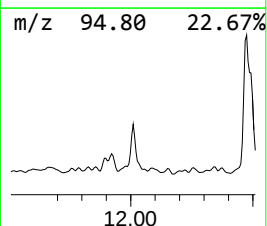
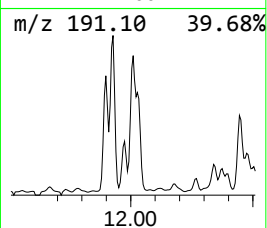
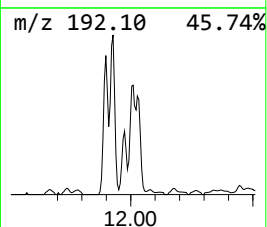
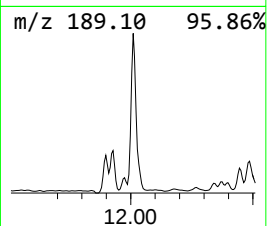
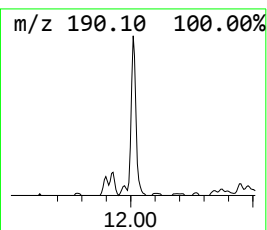
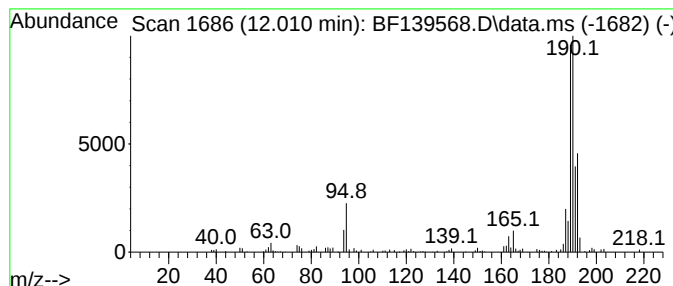
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.010	3.58 ng	133991	Phenanthrene-d10	11.398	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	86
2	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	43
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	37
4	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	32
5	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

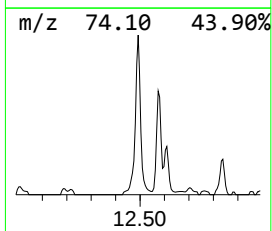
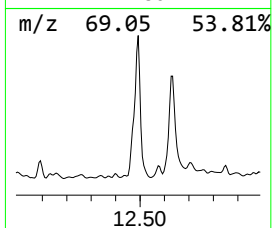
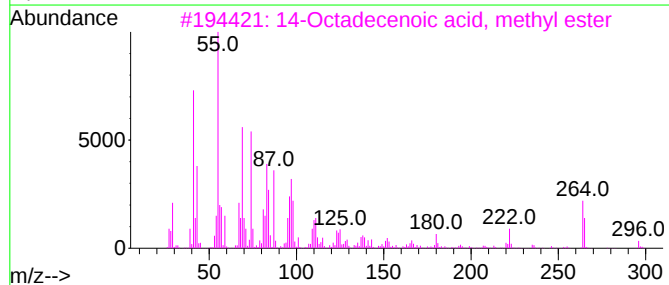
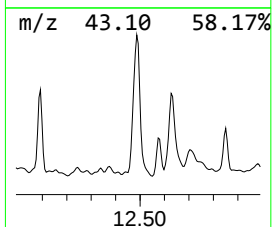
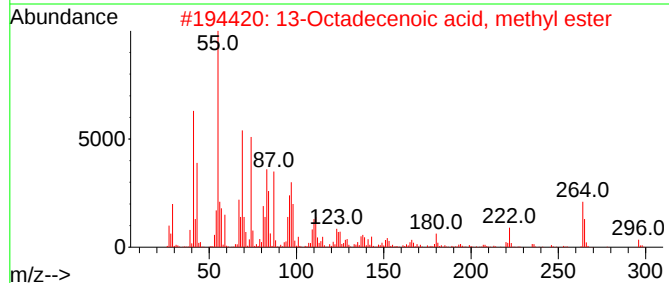
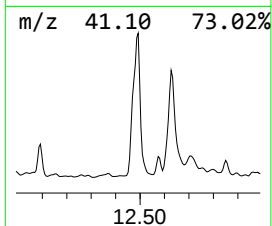
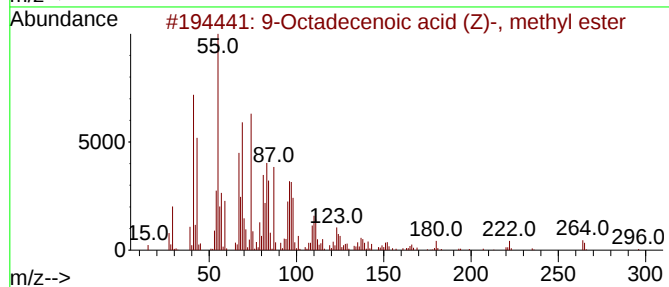
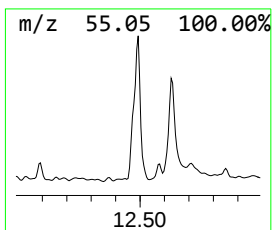
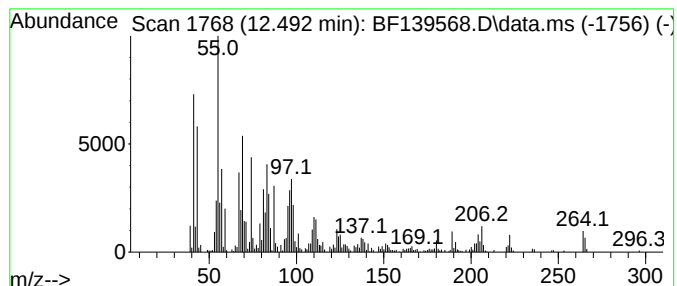
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.492	17.76 ng	664961	Phenanthrene-d10	11.398	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	98
3	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	97
4	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	96
5	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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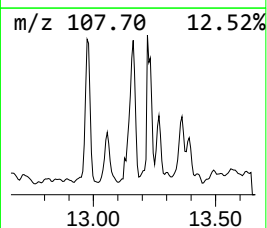
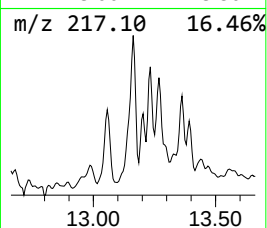
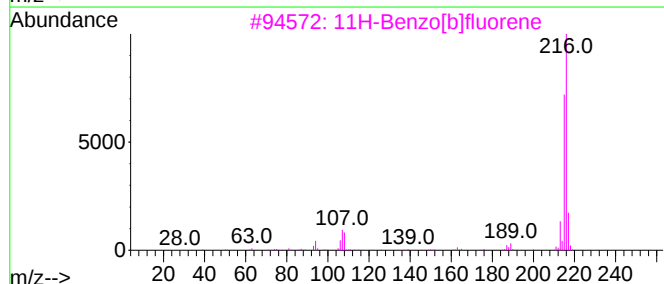
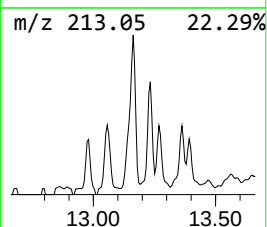
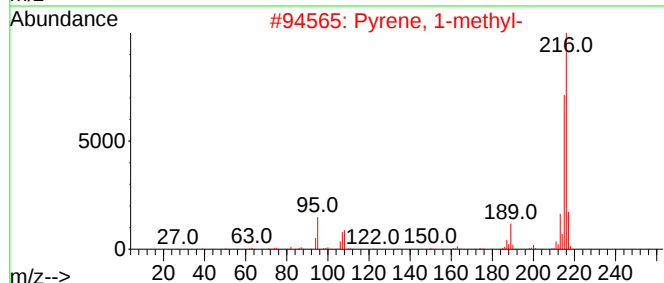
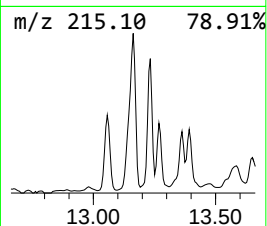
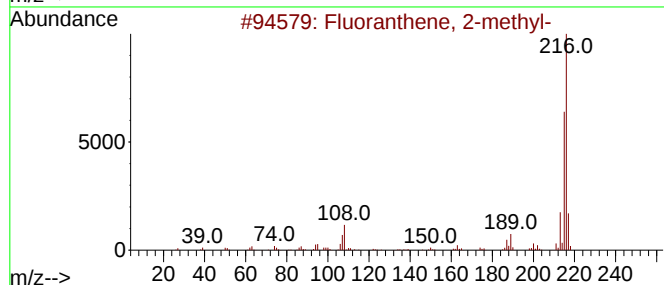
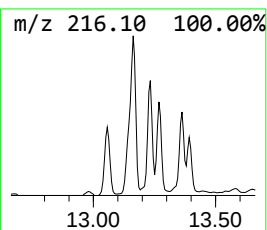
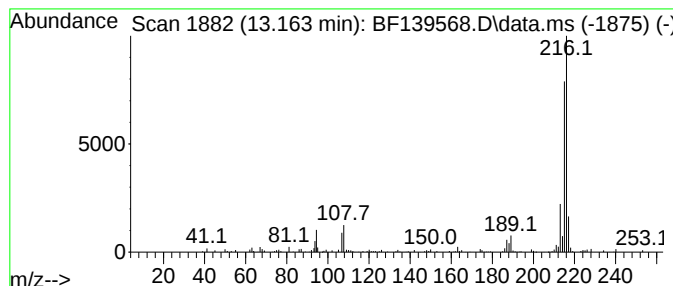
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	2.31 ng	110641	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-092524

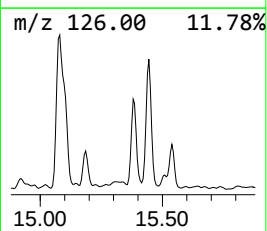
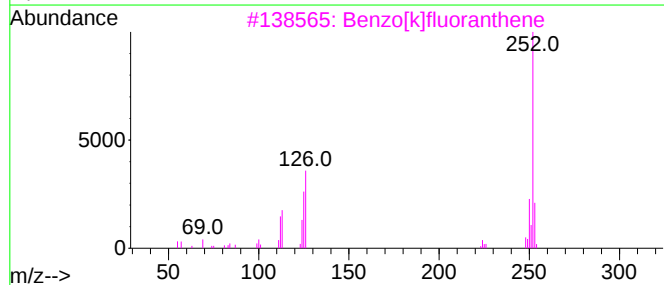
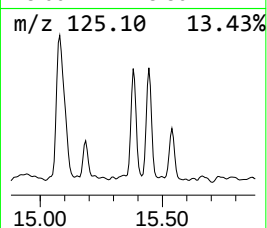
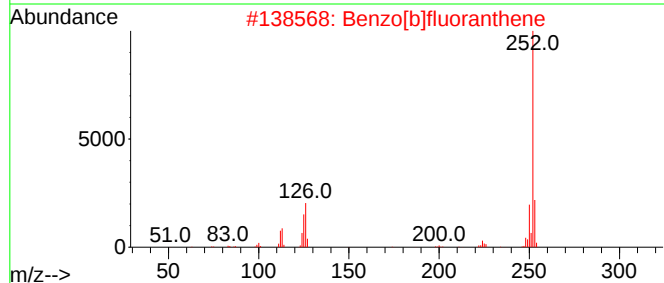
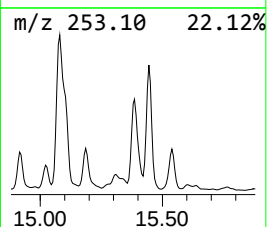
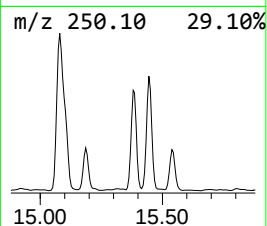
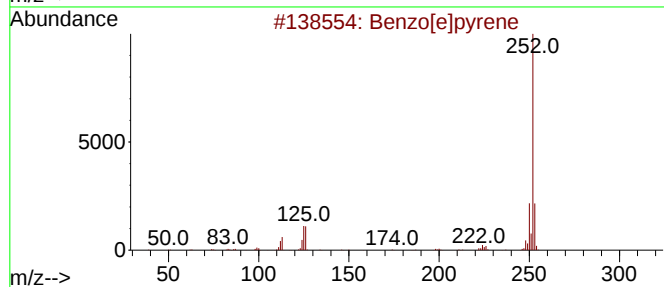
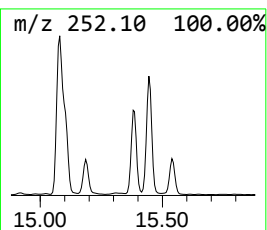
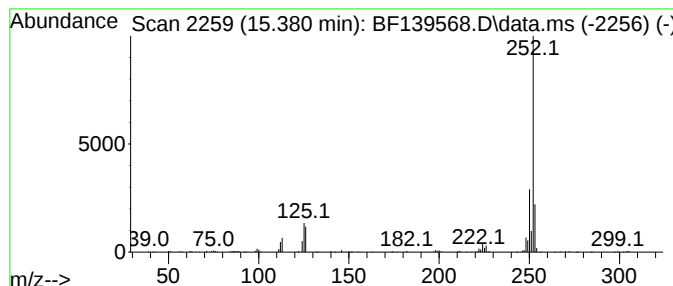
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.380	6.58 ng	153759	Perylene-d12	15.509

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	96
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
5			Benzo[a]pyrene	252	C20H12	000050-32-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092624\
Data File : BF139568.D
Acq On : 27 Sep 2024 01:42
Operator : RC/JU
Sample : P4197-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-092524

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...	2.134	24.3	ng	639234	1	6.875	525215	20.0
Dodecane	10.333	3.3	ng	113361	3	9.910	689747	20.0
Tridecane	10.810	4.6	ng	171635	4	11.398	748832	20.0
Nonadecane	11.257	2.7	ng	102660	4	11.398	748832	20.0
Tridecane, 1-iodo-	11.686	2.2	ng	82273	4	11.398	748832	20.0
Hexadecanoic ac...	11.786	2.8	ng	103656	4	11.398	748832	20.0
Phenanthrene, 1...	11.921	4.2	ng	155757	4	11.398	748832	20.0
4H-Cyclopenta[d...	12.010	3.6	ng	133991	4	11.398	748832	20.0
9-Octadecenoic ...	12.492	17.8	ng	664961	4	11.398	748832	20.0
Fluoranthene, 2...	13.163	2.3	ng	110641	5	14.033	958326	20.0
Benzo[e]pyrene	15.380	6.6	ng	153759	6	15.509	467499	20.0