

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
 Data File : BF128419.D
 Acq On : 24 May 2022 00:37
 Operator : CG\JU
 Sample : N2964-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-051922

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-BF052022.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128419.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.098	475	478	489	rVB	28129	39433	1.56%	0.160%
2	5.481	540	543	563	rBV	2192610	2381122	94.45%	9.638%
3	6.492	711	715	727	rBV	2465528	2387445	94.70%	9.664%
4	6.675	743	746	749	rBV	38324	33743	1.34%	0.137%
5	6.857	773	777	783	rBV	775625	709856	28.16%	2.873%
6	7.116	816	821	824	rBV2	45932	47162	1.87%	0.191%
7	7.381	861	866	869	rBV2	20798	29137	1.16%	0.118%
8	7.422	869	873	880	rBV	1786396	1699963	67.43%	6.881%
9	7.804	935	938	941	rVB2	24623	31777	1.26%	0.129%
10	7.875	947	950	952	rBV	55290	53403	2.12%	0.216%
11	8.039	974	978	986	rBV	443516	521901	20.70%	2.112%
12	8.104	986	989	991	rVV	253851	222251	8.82%	0.900%
13	8.139	991	995	997	rVV	1035002	941805	37.36%	3.812%
14	8.157	997	998	1001	rVV	227317	193731	7.68%	0.784%
15	8.181	1001	1002	1005	rVB	103393	70342	2.79%	0.285%
16	8.216	1005	1008	1010	rBV3	30918	32648	1.30%	0.132%
17	8.334	1024	1028	1034	rVB2	33625	35872	1.42%	0.145%
18	8.416	1039	1042	1047	rVB4	60872	78779	3.12%	0.319%
19	8.469	1047	1051	1053	rBV3	26105	31841	1.26%	0.129%
20	8.498	1054	1056	1061	rVV2	36276	43308	1.72%	0.175%
21	8.539	1061	1063	1066	rVB	99023	82826	3.29%	0.335%
22	8.639	1076	1080	1082	rBV2	52240	46999	1.86%	0.190%
23	8.716	1090	1093	1097	rVB	270335	218328	8.66%	0.884%
24	8.775	1097	1103	1105	rBV6	17049	36161	1.43%	0.146%
25	8.810	1105	1109	1113	rVB3	43821	68277	2.71%	0.276%
26	8.851	1113	1116	1119	rBV	223661	173121	6.87%	0.701%
27	8.951	1128	1133	1138	rBV	208295	182017	7.22%	0.737%
28	8.998	1138	1141	1142	rVV2	33401	29694	1.18%	0.120%
29	9.039	1145	1148	1151	rVB3	29482	43224	1.71%	0.175%
30	9.081	1151	1155	1157	rBV2	50016	47888	1.90%	0.194%
31	9.145	1163	1166	1169	rVV	84107	58586	2.32%	0.237%
32	9.210	1173	1177	1181	rBV	2583456	2521067	100.00%	10.204%
33	9.281	1185	1189	1192	rBV	262954	207043	8.21%	0.838%
34	9.322	1193	1196	1200	rVV	48988	55466	2.20%	0.225%
35	9.357	1200	1202	1205	rVB3	34428	27523	1.09%	0.111%
36	9.481	1219	1223	1227	rBV3	69684	95862	3.80%	0.388%

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37	9.551	1227	1235	1238	rVV3	100930	147482	5.85%	0.597%
38	9.581	1238	1240	1241	rVV	67772	60024	2.38%	0.243%
39	9.610	1241	1245	1251	rVV2	138471	205832	8.16%	0.833%
40	9.669	1251	1255	1260	rVB2	46920	58283	2.31%	0.236%

41	9.757	1266	1270	1274	rBV	99589	91653	3.64%	0.371%
42	9.810	1276	1279	1283	rVB	165335	125670	4.98%	0.509%
43	9.892	1290	1293	1296	rVV	1037015	892511	35.40%	3.613%
44	9.928	1296	1299	1306	rVB	150944	143024	5.67%	0.579%
45	9.992	1306	1310	1315	rVB5	15687	29332	1.16%	0.119%

46	10.051	1315	1320	1321	rBV3	20402	28950	1.15%	0.117%
47	10.104	1325	1329	1331	rVB2	71370	69045	2.74%	0.279%
48	10.133	1331	1334	1338	rVB3	38832	40822	1.62%	0.165%
49	10.210	1343	1347	1350	rBV2	29500	36304	1.44%	0.147%
50	10.310	1358	1364	1368	rBV	105643	104936	4.16%	0.425%

51	10.445	1384	1387	1393	rVB	185523	164348	6.52%	0.665%
52	10.528	1396	1401	1404	rVB3	40776	53461	2.12%	0.216%
53	10.686	1425	1428	1442	rVB	1393040	1313815	52.11%	5.318%
54	10.786	1442	1445	1451	rBV	53080	62380	2.47%	0.252%
55	10.992	1475	1480	1483	rBV2	33327	46943	1.86%	0.190%

56	11.233	1518	1521	1524	rBV	43420	39628	1.57%	0.160%
57	11.286	1527	1530	1534	rVB	53184	46042	1.83%	0.186%
58	11.386	1544	1547	1549	rBV	957557	803367	31.87%	3.252%
59	11.410	1549	1551	1555	rVV	607076	517034	20.51%	2.093%
60	11.463	1555	1560	1565	rVV	155292	163959	6.50%	0.664%

61	11.627	1583	1588	1591	rBV	32624	35676	1.42%	0.144%
62	11.763	1608	1611	1614	rBV	86902	64111	2.54%	0.260%
63	11.892	1629	1633	1635	rBV	65030	65723	2.61%	0.266%
64	11.922	1635	1638	1642	rVV	75174	72078	2.86%	0.292%
65	11.969	1643	1646	1649	rVV	34088	31610	1.25%	0.128%

66	12.004	1649	1652	1654	rVV	126920	125389	4.97%	0.508%
67	12.027	1654	1656	1660	rVB	41973	37922	1.50%	0.153%
68	12.186	1680	1683	1686	rBV	38457	32941	1.31%	0.133%
69	12.451	1723	1728	1730	rBV3	96970	121600	4.82%	0.492%
70	12.474	1730	1732	1735	rVB2	89197	79741	3.16%	0.323%

71	12.557	1744	1746	1750	rVB2	29275	28398	1.13%	0.115%
72	12.604	1751	1754	1759	rBV	483172	410989	16.30%	1.664%
73	12.780	1782	1784	1787	rVB2	31148	25778	1.02%	0.104%
74	12.833	1787	1793	1799	rBV	371926	446440	17.71%	1.807%
75	12.974	1812	1817	1820	rBV	1824895	1726591	68.49%	6.989%

76	13.051	1826	1830	1835	rBV3	22711	39881	1.58%	0.161%
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77	13.163	1843	1849	1852	rBV	70301	86463	3.43%	0.350%
78	13.227	1858	1860	1863	rBV	57367	51615	2.05%	0.209%
79	13.269	1864	1867	1869	rBV2	27728	27339	1.08%	0.111%
80	13.392	1886	1888	1893	rVB2	21294	25350	1.01%	0.103%
81	13.527	1908	1911	1914	rBV2	27954	31151	1.24%	0.126%
82	13.857	1962	1967	1970	rVB3	39729	51372	2.04%	0.208%
83	13.910	1971	1976	1979	rBV6	32366	62059	2.46%	0.251%
84	14.039	1993	1998	2000	rBV2	683300	778836	30.89%	3.152%
85	14.063	2000	2002	2009	rVB	158172	147808	5.86%	0.598%
86	14.127	2010	2013	2014	rBV2	32940	37050	1.47%	0.150%
87	14.174	2018	2021	2023	rBV	36116	30211	1.20%	0.122%
88	14.786	2122	2125	2128	rVB	40468	37212	1.48%	0.151%
89	15.086	2173	2176	2180	rBV	120313	174597	6.93%	0.707%
90	15.121	2180	2182	2187	rVB2	88388	92347	3.66%	0.374%
91	15.204	2193	2196	2200	rBV	34025	42480	1.69%	0.172%
92	15.398	2226	2229	2234	rVB	71164	81154	3.22%	0.328%
93	15.462	2237	2240	2244	rBV	104871	119417	4.74%	0.483%
94	15.521	2245	2250	2255	rBV2	468045	609421	24.17%	2.467%
95	17.045	2504	2509	2520	rVB3	47489	99736	3.96%	0.404%
96	17.498	2583	2586	2592	rVB2	27089	50626	2.01%	0.205%

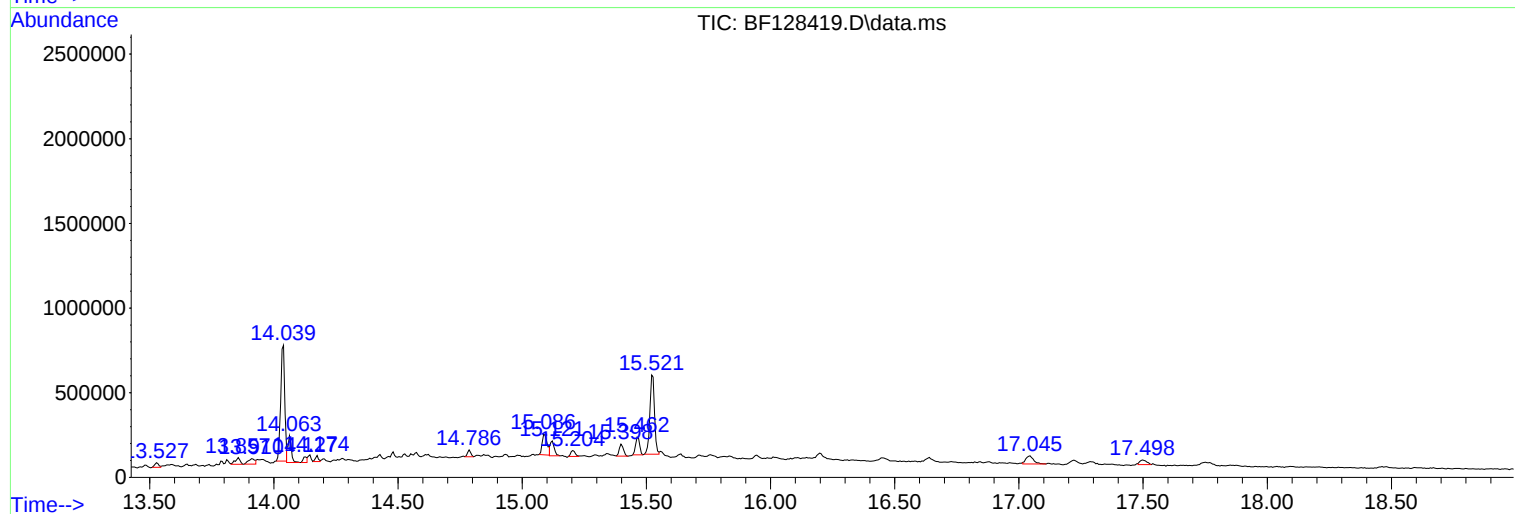
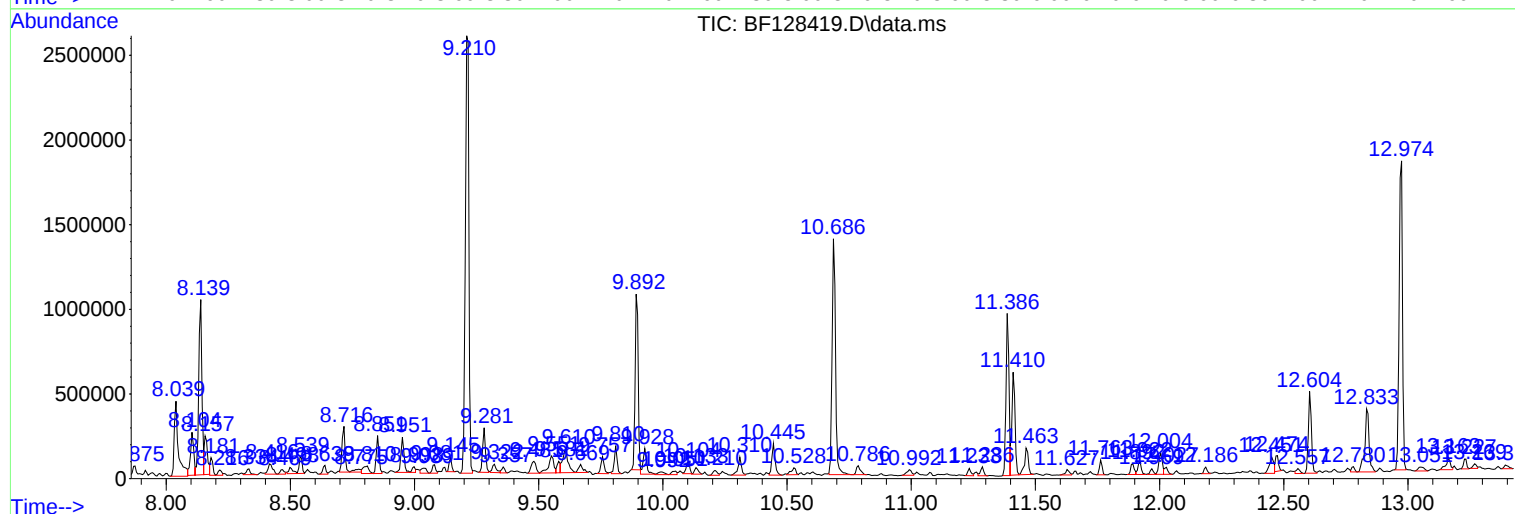
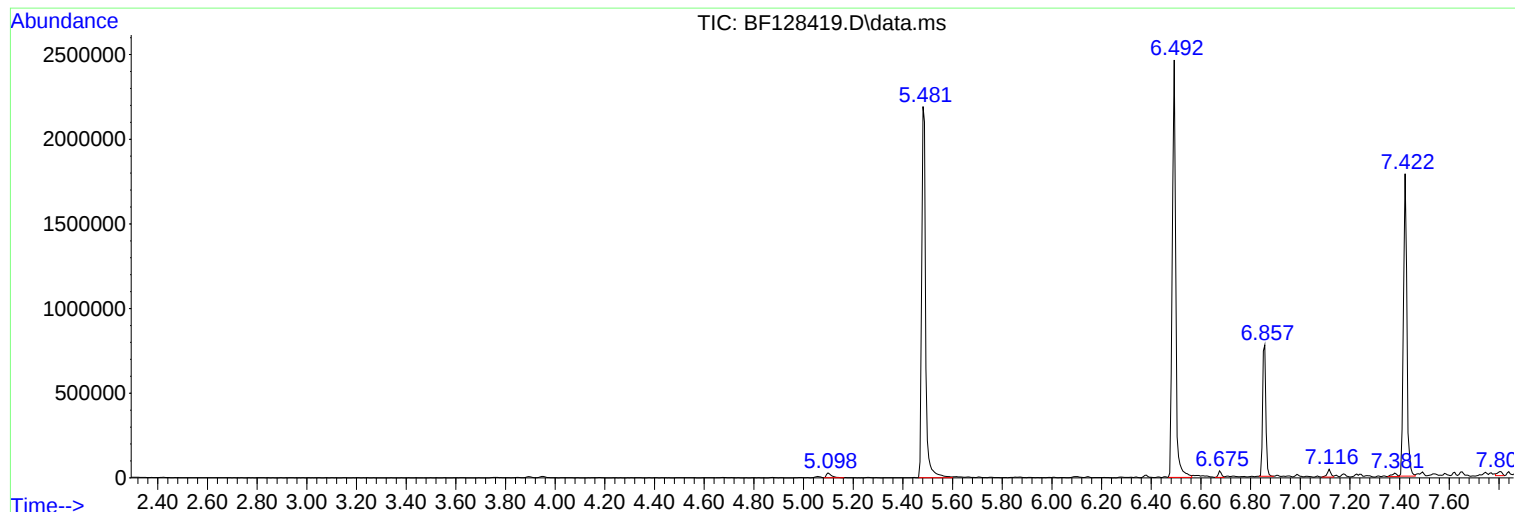
Sum of corrected areas: 24705558

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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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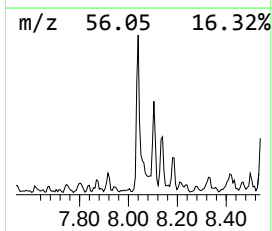
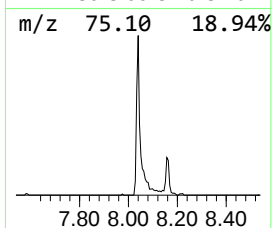
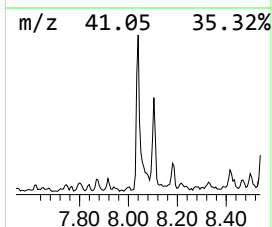
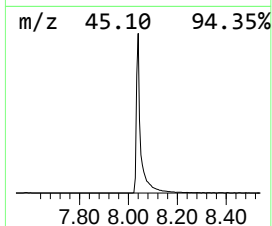
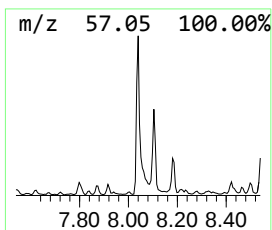
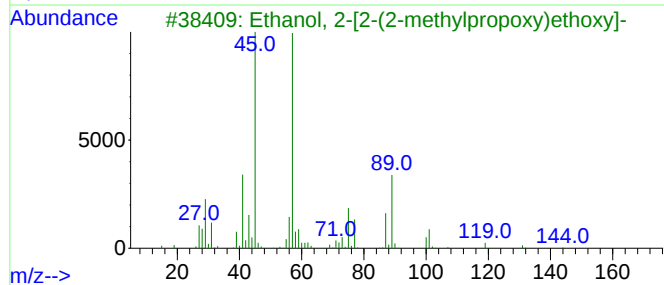
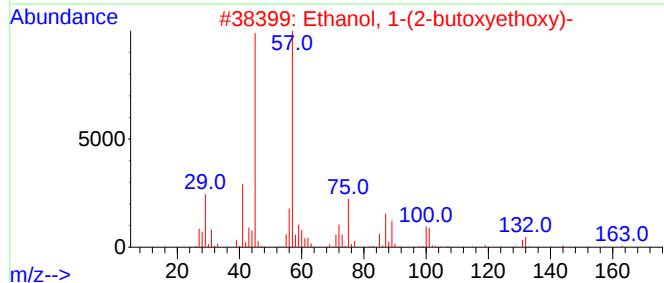
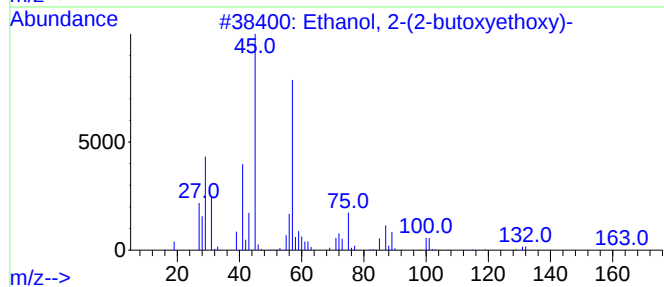
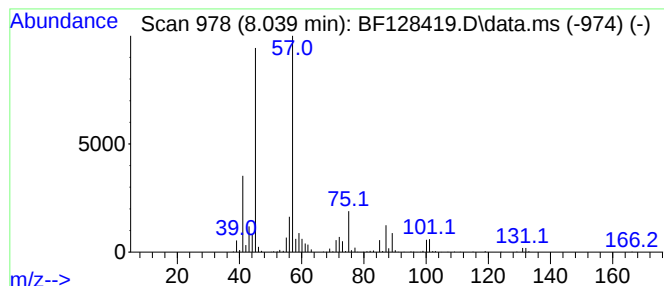
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.039	11.08 ng	521901	Naphthalene-d8	8.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	59



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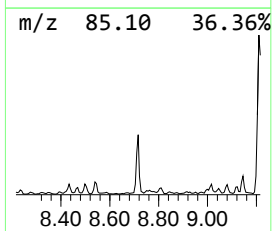
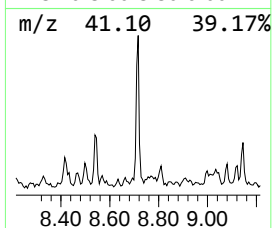
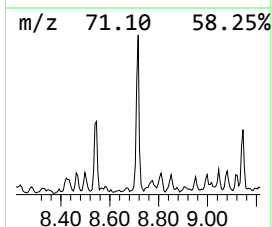
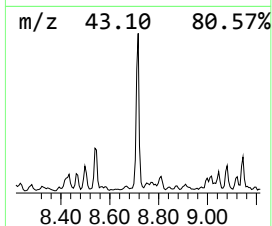
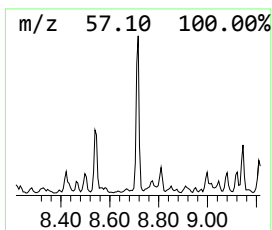
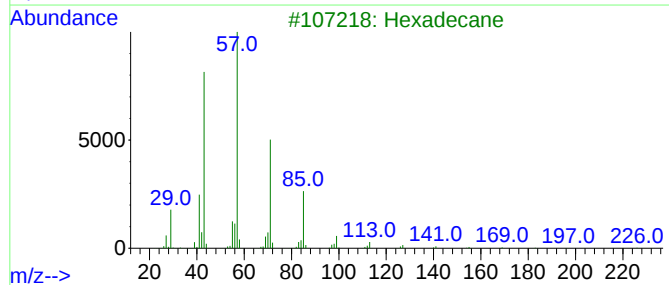
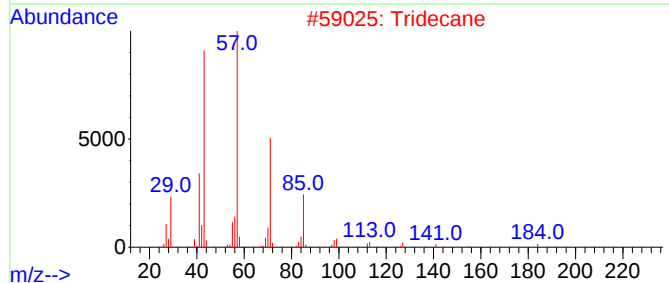
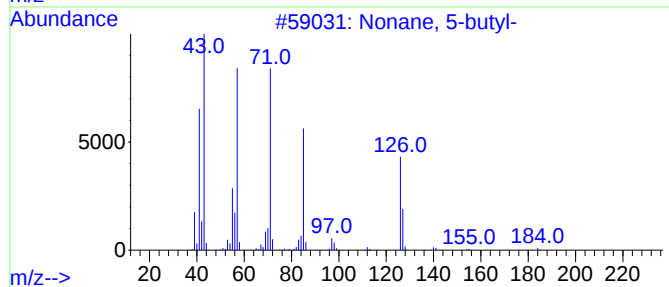
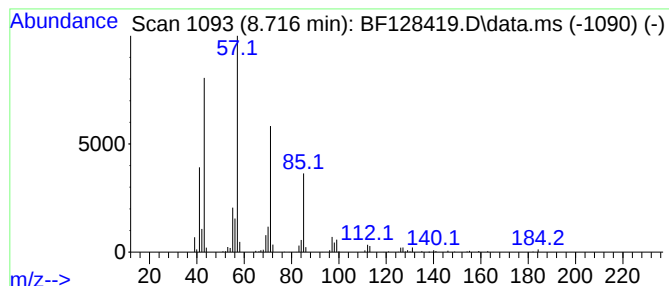
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Nonane, 5-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.716	4.64 ng	218328	Naphthalene-d8	8.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 5-butyl-	184	C13H28	017312-63-9	90
2			Tridecane	184	C13H28	000629-50-5	87
3			Hexadecane	226	C16H34	000544-76-3	87
4			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	86
5			Tetradecane	198	C14H30	000629-59-4	80



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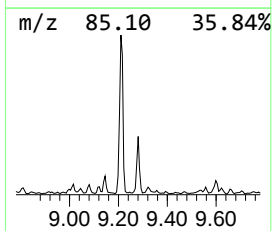
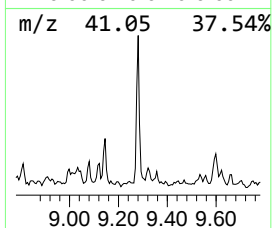
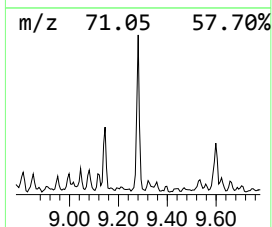
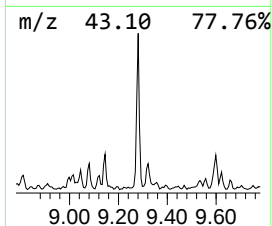
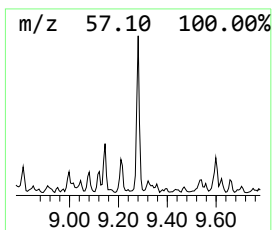
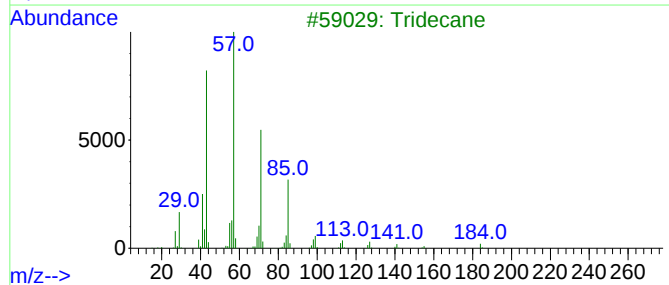
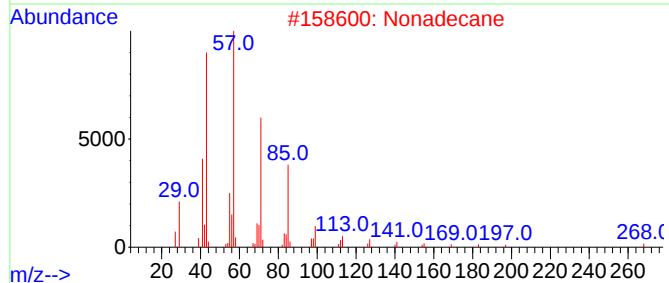
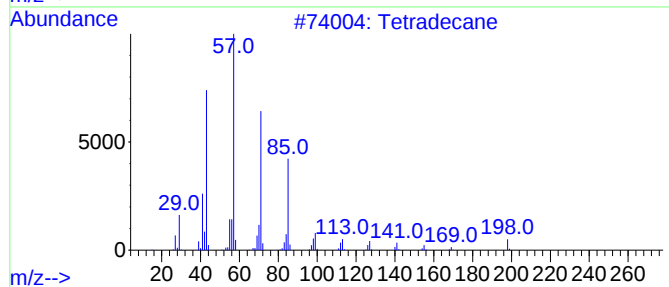
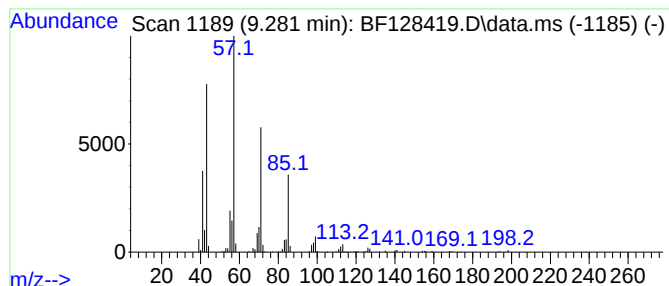
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tetradecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.281	4.64 ng	207043	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	94
2			Nonadecane	268	C19H40	000629-92-5	91
3			Tridecane	184	C13H28	000629-50-5	90
4			Hexadecane	226	C16H34	000544-76-3	90
5			Nonane, 5-butyl-	184	C13H28	017312-63-9	80



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Data File : BF128419.D
Acq On : 24 May 2022 00:37
Operator : CG\JU
Sample : N2964-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-01-051922

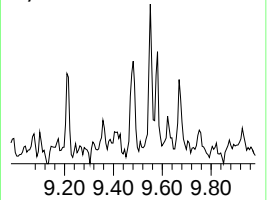
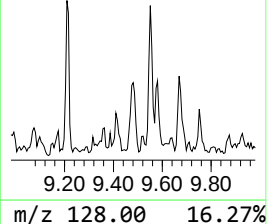
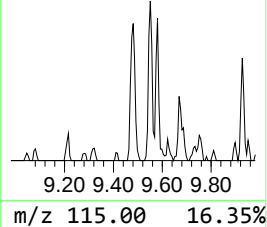
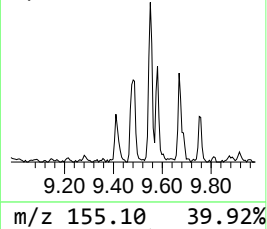
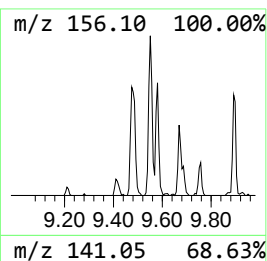
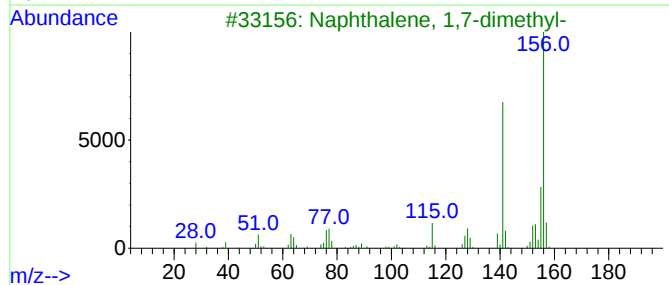
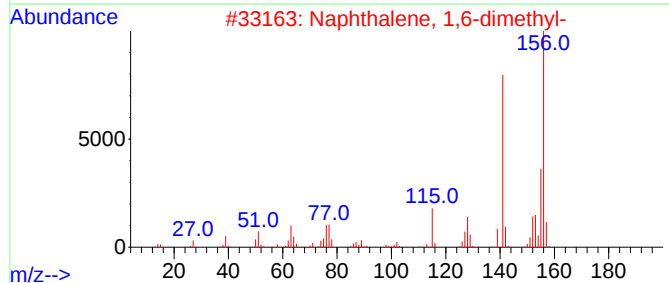
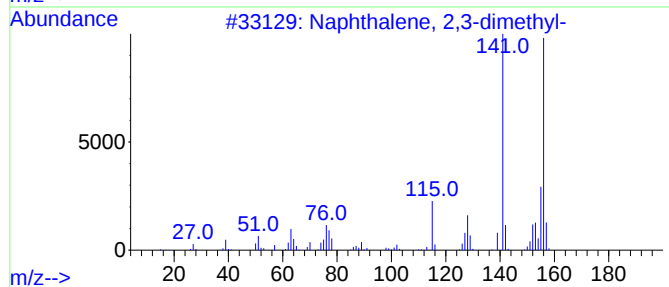
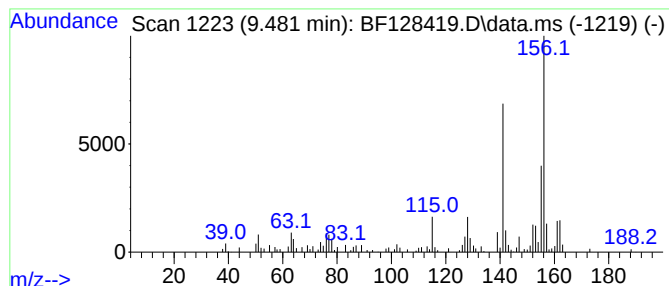
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.481	2.15 ng	95862	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
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ClientSampleId :
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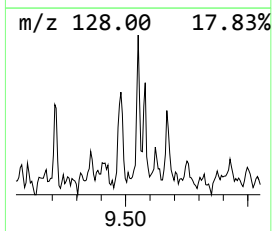
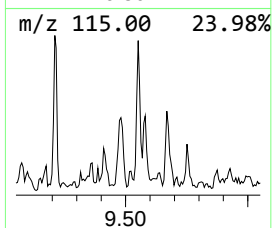
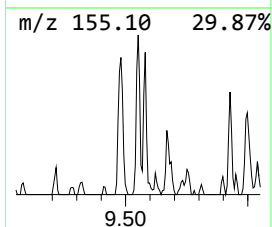
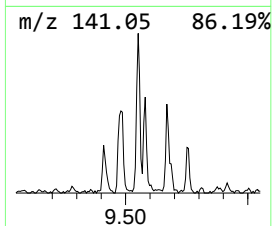
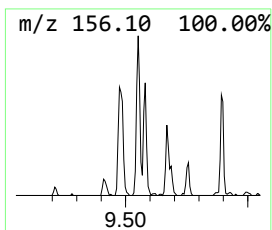
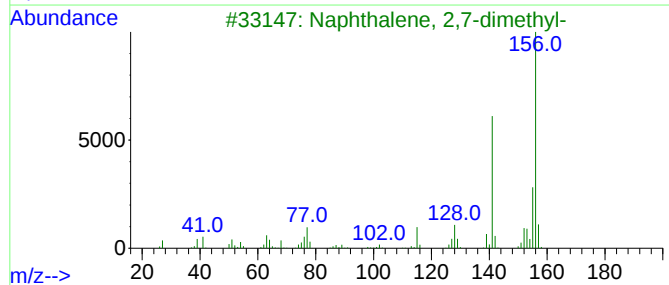
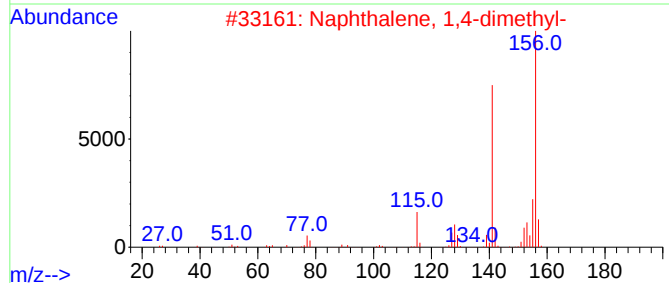
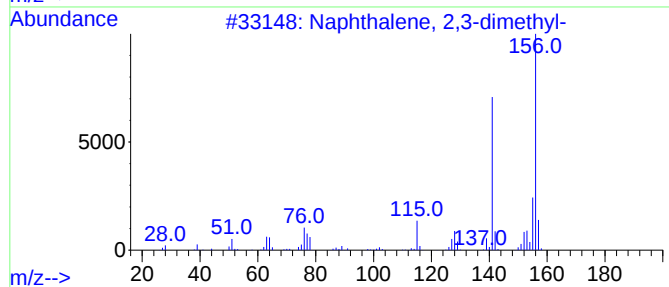
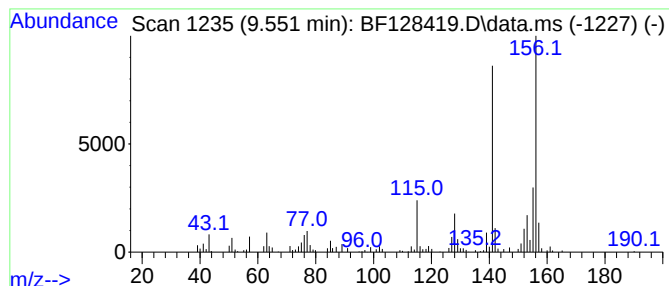
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.551	3.30 ng	147482	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
Acq On : 24 May 2022 00:37
Operator : CG\JU
Sample : N2964-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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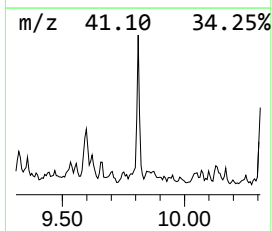
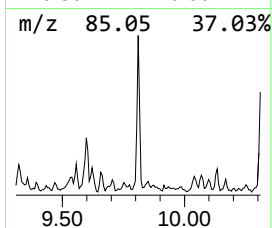
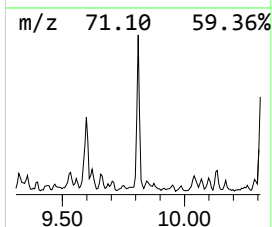
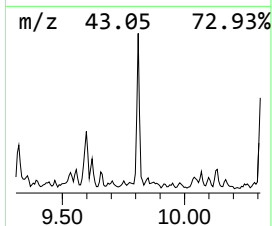
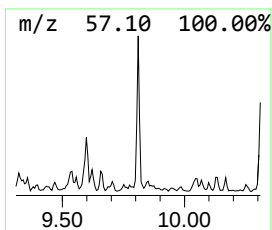
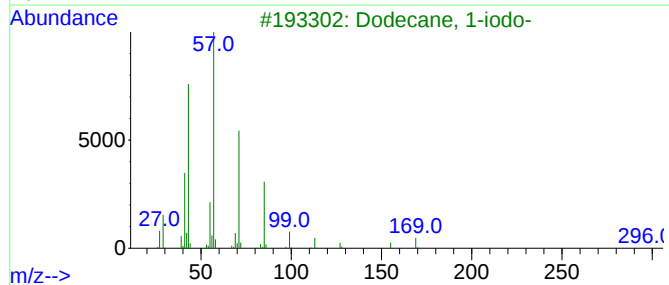
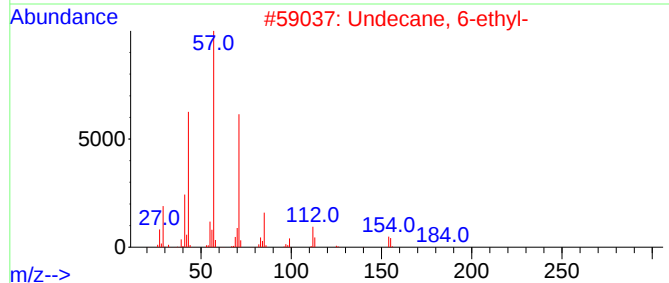
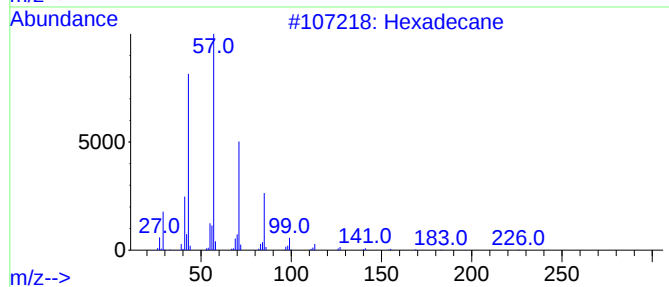
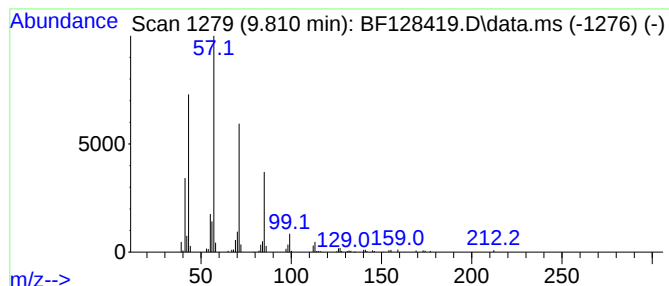
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.810	2.82 ng	125670	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	90
2			Undecane, 6-ethyl-	184	C13H28	017312-60-6	72
3			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	64
4			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
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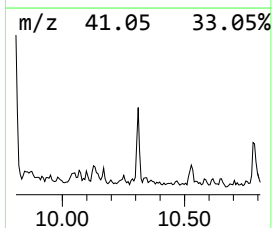
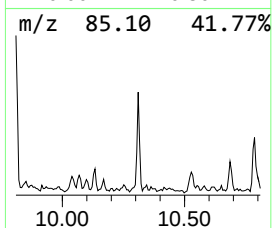
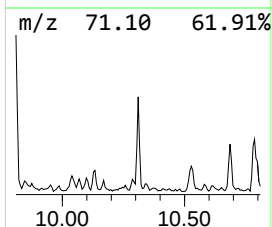
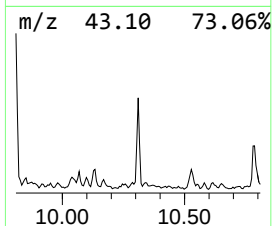
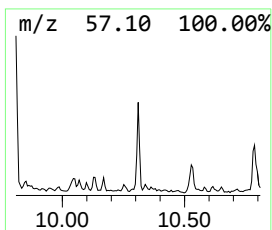
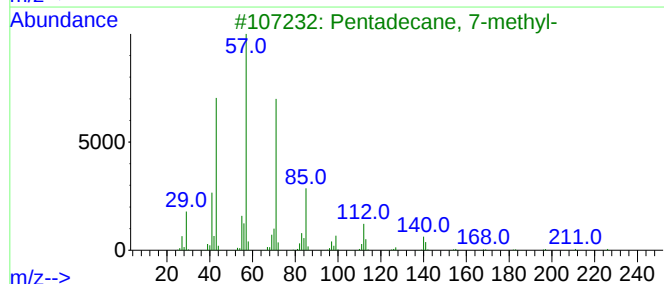
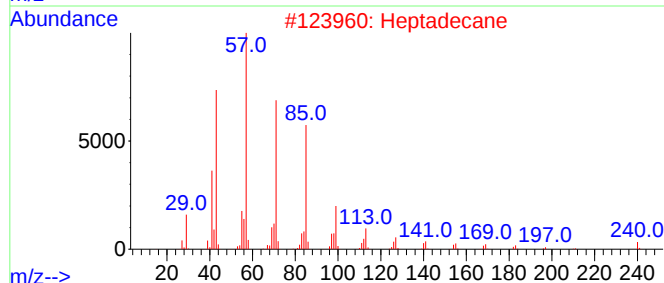
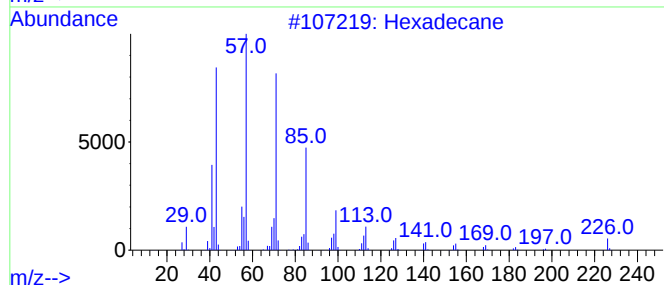
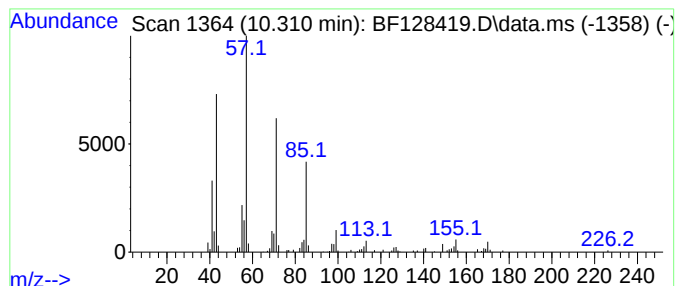
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.310	2.35 ng	104936	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Heptadecane	240	C17H36	000629-78-7	86
3			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	76
4			Tridecane, 7-propyl-	226	C16H34	055045-09-5	72
5			Decane, 3-methyl-	156	C11H24	013151-34-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
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ClientSampleId :
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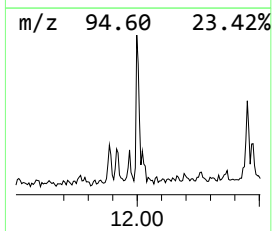
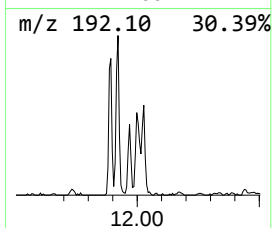
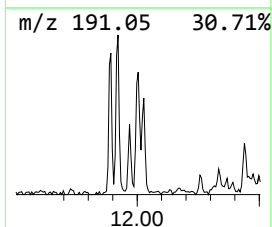
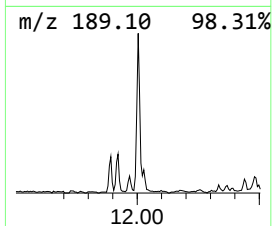
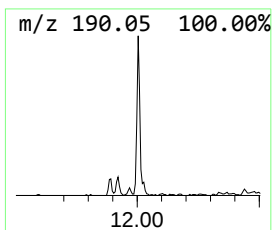
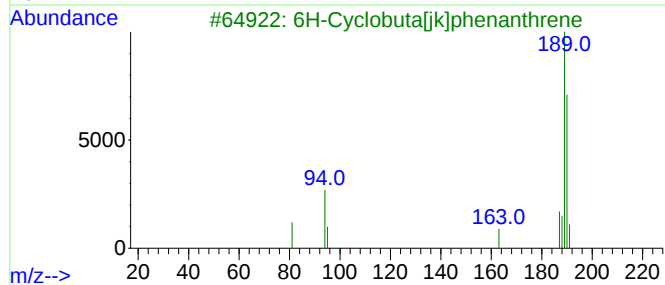
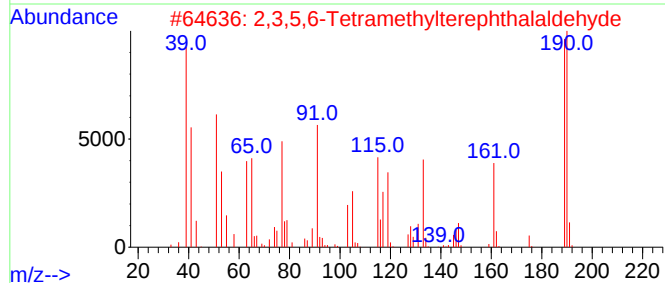
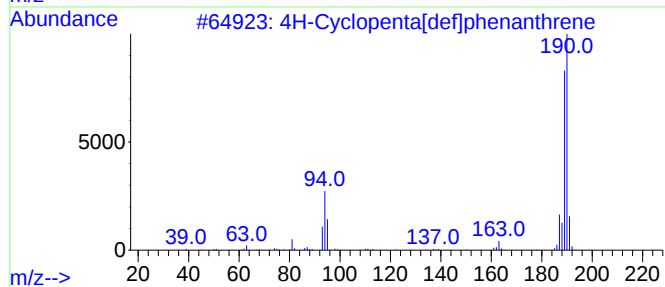
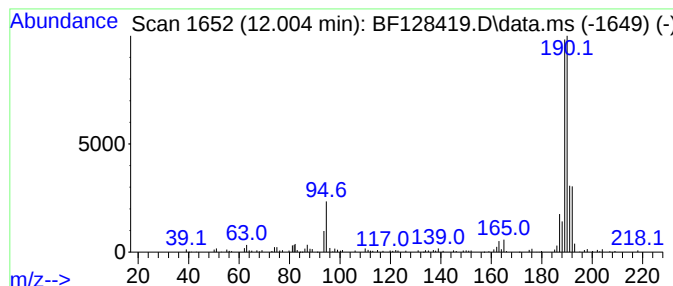
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	3.12 ng	125389	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	52
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
4			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
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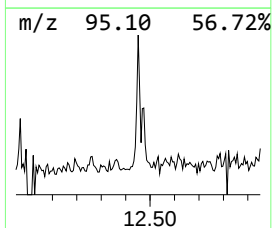
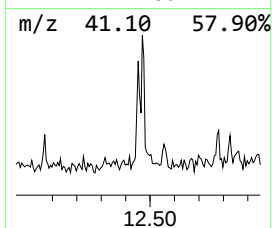
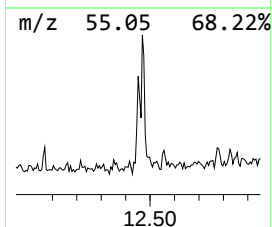
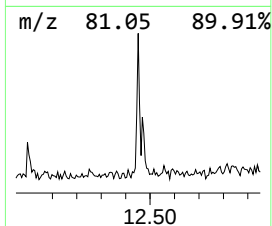
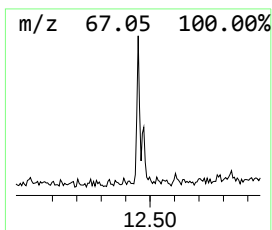
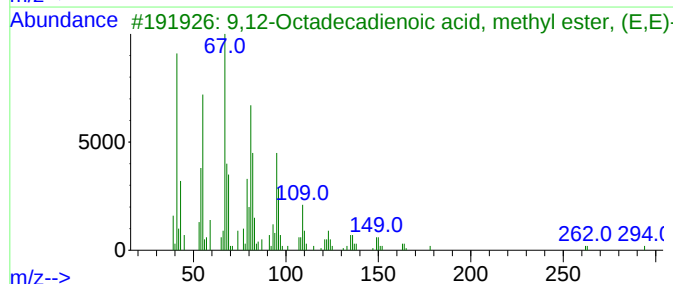
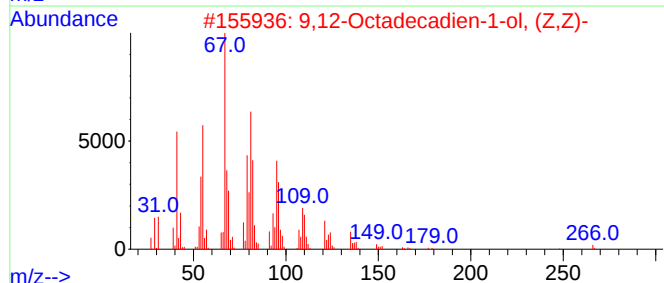
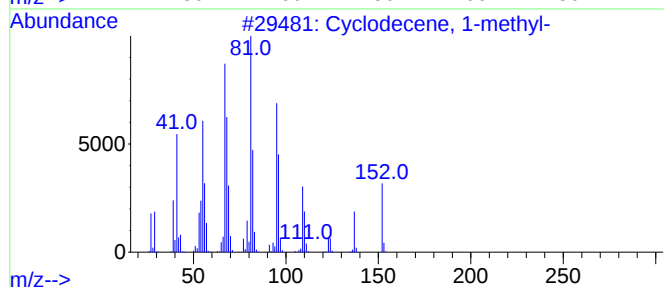
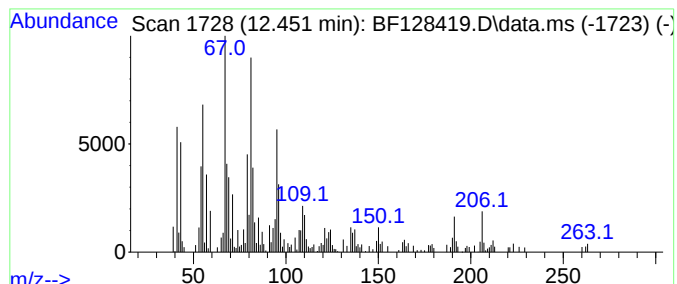
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclodecene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.451	3.03 ng	121600	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	70
2			9,12-Octadecadien-1-ol, (Z,Z)-	266	C18H34O	000506-43-4	64
3			9,12-Octadecadienoic acid, methyl...	294	C19H34O2	002566-97-4	62
4			4-Methyl-exo-tricyclo[6.2.1.0(2...	164	C12H20	1000215-29-8	58
5			11,14-Eicosadienoic acid, methyl...	322	C21H38O2	002463-02-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
Acq On : 24 May 2022 00:37
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ALS Vial : 17 Sample Multiplier: 1

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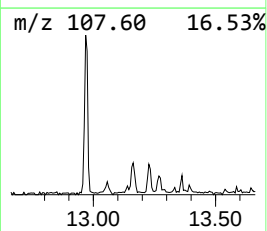
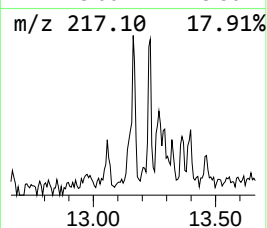
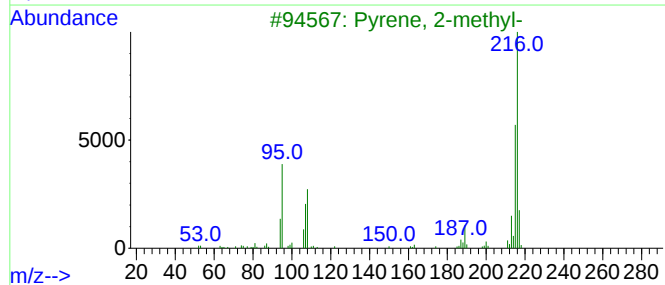
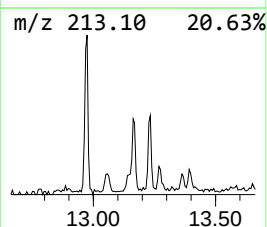
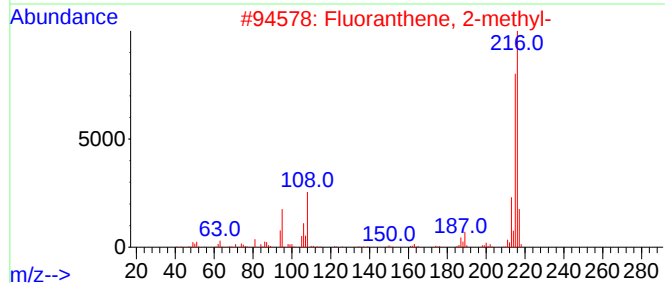
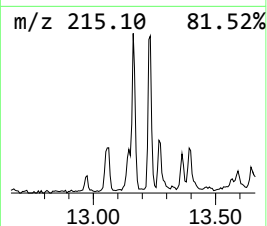
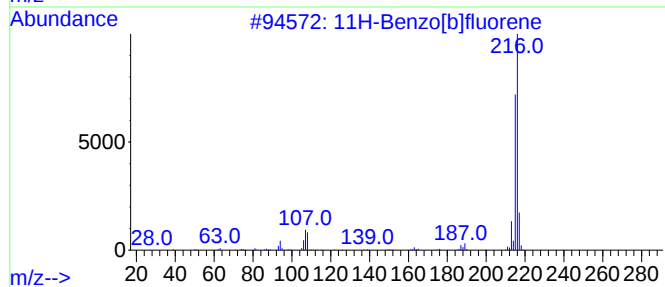
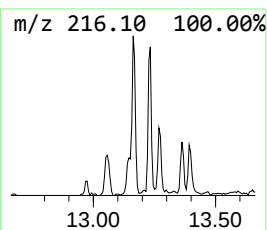
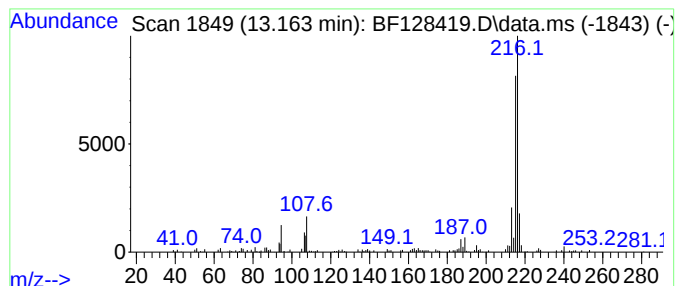
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	2.22 ng	86463	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
Acq On : 24 May 2022 00:37
Operator : CG\JU
Sample : N2964-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-051922

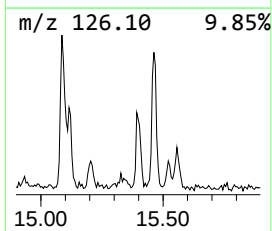
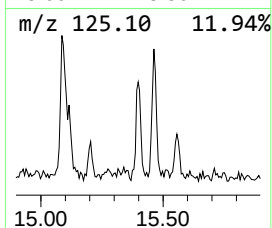
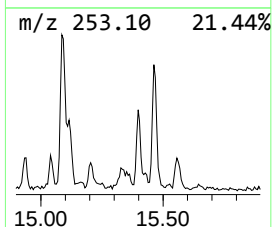
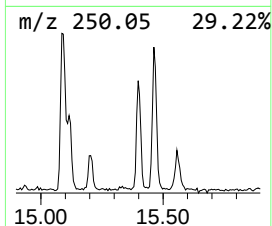
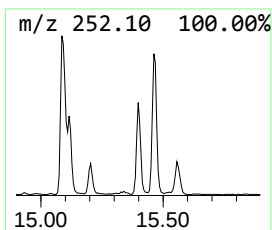
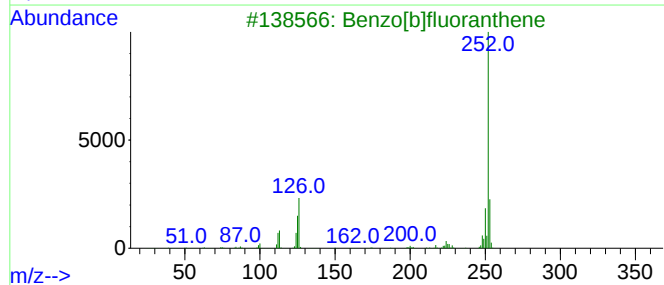
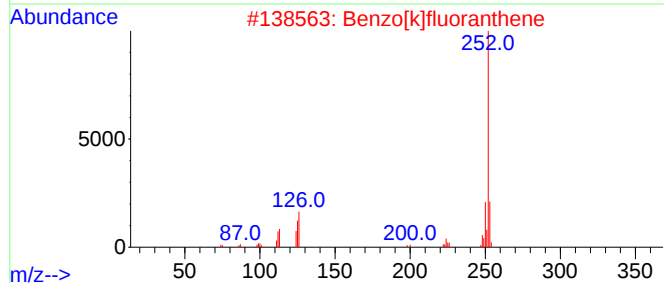
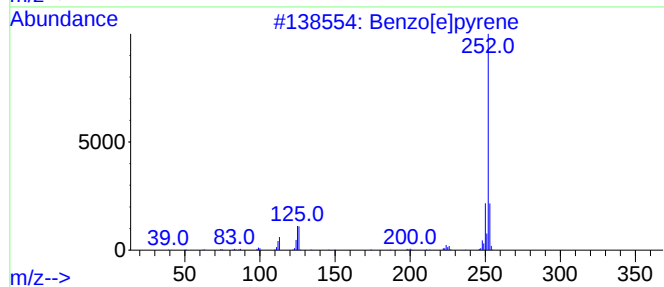
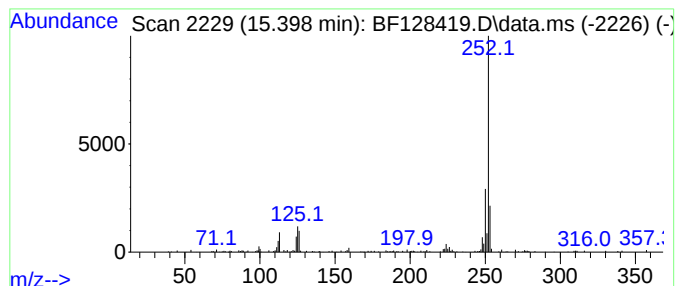
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.398	2.66 ng	81154	Perylene-d12	15.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
4			Benzo[a]pyrene	252	C20H12	000050-32-8	90
5			Perylene	252	C20H12	000198-55-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052322\
Data File : BF128419.D
Acq On : 24 May 2022 00:37
Operator : CG\JU
Sample : N2964-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-051922

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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
Ethanol, 2-(2-b...	8.039	11.1	ng	521901	2	8.139	941805	20.0
Nonane, 5-butyl-	8.716	4.6	ng	218328	2	8.139	941805	20.0
Tetradecane	9.281	4.6	ng	207043	3	9.892	892511	20.0
Naphthalene, 2,...	9.481	2.1	ng	95862	3	9.892	892511	20.0
Naphthalene, 1,...	9.551	3.3	ng	147482	3	9.892	892511	20.0
Hexadecane	9.810	2.8	ng	125670	3	9.892	892511	20.0
Heptadecane	10.310	2.4	ng	104936	3	9.892	892511	20.0
4H-Cyclopenta[d...]	12.004	3.1	ng	125389	4	11.386	803367	20.0
Cyclodecene, 1-...	12.451	3.0	ng	121600	4	11.386	803367	20.0
11H-Benzo[b]flu...	13.163	2.2	ng	86463	5	14.039	778836	20.0
Benzo[e]pyrene	15.398	2.7	ng	81154	6	15.527	609421	20.0