

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101477.D
 Acq On : 18 Dec 2017 13:42
 Operator : SJ/JU
 Sample : I6679-07 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-05-121317-C

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.010	494	497	512	rBV	65889	108870	8.32%	0.560%
2	5.422	564	567	587	rBV	482483	733924	56.12%	3.773%
3	6.445	738	741	760	rBV	511026	815234	62.34%	4.191%
4	6.581	760	764	770	rVV	753705	791283	60.51%	4.068%
5	6.628	770	772	778	rVB3	21894	30351	2.32%	0.156%
6	6.804	799	802	810	rBV	997999	936162	71.58%	4.813%
7	6.963	825	829	838	rVB	597513	598907	45.80%	3.079%
8	7.375	896	899	909	rBV	371930	494996	37.85%	2.545%
9	7.604	935	938	941	rBV3	16550	15580	1.19%	0.080%
10	7.828	973	976	978	rBV2	25529	23378	1.79%	0.120%
11	8.063	1013	1016	1017	rBV	52144	42632	3.26%	0.219%
12	8.087	1017	1020	1026	rVV	1150123	1214507	92.87%	6.244%
13	8.134	1026	1028	1032	rVB2	44970	54047	4.13%	0.278%
14	8.287	1051	1054	1056	rVB2	18371	14797	1.13%	0.076%
15	8.375	1064	1069	1073	rVB6	26582	33068	2.53%	0.170%
16	8.498	1087	1090	1092	rBV	48730	34701	2.65%	0.178%
17	8.592	1103	1106	1108	rBV2	30197	31274	2.39%	0.161%
18	8.669	1116	1119	1123	rBV	87380	67912	5.19%	0.349%
19	8.751	1128	1133	1139	rVB5	25342	53137	4.06%	0.273%
20	8.910	1157	1160	1164	rBV2	42429	47392	3.62%	0.244%
21	8.986	1170	1173	1177	rVB4	16259	20802	1.59%	0.107%
22	9.034	1177	1181	1185	rBV4	24899	29101	2.23%	0.150%
23	9.098	1190	1192	1195	rVB	56419	39825	3.05%	0.205%
24	9.163	1199	1203	1211	rBV	963882	921244	70.44%	4.736%
25	9.234	1212	1215	1220	rVV	107234	92417	7.07%	0.475%
26	9.310	1226	1228	1231	rVB3	24521	19151	1.46%	0.098%
27	9.439	1245	1250	1253	rBV4	21007	30968	2.37%	0.159%
28	9.504	1257	1261	1264	rVV3	37763	51505	3.94%	0.265%
29	9.557	1264	1270	1277	rVB4	67547	150270	11.49%	0.773%
30	9.622	1277	1281	1283	rBV5	18511	18176	1.39%	0.093%
31	9.710	1292	1296	1300	rBV	84521	93322	7.14%	0.480%
32	9.763	1303	1305	1308	rVB	88900	66357	5.07%	0.341%
33	9.845	1315	1319	1323	rBV	1435026	1302308	99.58%	6.695%
34	9.875	1323	1324	1332	rVB2	58467	68981	5.27%	0.355%

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35	9.945	1333	1336	1340	rVB4	15666	24132	1.85%	0.124%
36	9.992	1340	1344	1347	rBV7	17715	27865	2.13%	0.143%
37	10.051	1351	1354	1357	rVB3	30426	33198	2.54%	0.171%
38	10.086	1357	1360	1364	rBV4	36758	36832	2.82%	0.189%
39	10.157	1369	1372	1376	rBV4	30235	39796	3.04%	0.205%
40	10.192	1376	1378	1383	rVB2	14816	21347	1.63%	0.110%
41	10.263	1383	1390	1393	rBV2	89269	100854	7.71%	0.518%
42	10.369	1406	1408	1410	rBV2	14615	14805	1.13%	0.076%
43	10.398	1410	1413	1419	rVB	45014	58496	4.47%	0.301%
44	10.481	1422	1427	1429	rBV2	46241	55552	4.25%	0.286%
45	10.557	1438	1440	1446	rVB6	23241	22951	1.75%	0.118%
46	10.639	1450	1454	1463	rBV	415717	457487	34.98%	2.352%
47	10.733	1467	1470	1477	rBV3	85282	132094	10.10%	0.679%
48	10.945	1500	1506	1509	rBV5	29204	52780	4.04%	0.271%
49	11.022	1517	1519	1522	rVB2	20322	18328	1.40%	0.094%
50	11.069	1522	1527	1530	rBV6	14617	31389	2.40%	0.161%
51	11.180	1544	1546	1549	rBV	73153	59047	4.52%	0.304%
52	11.210	1549	1551	1553	rVV2	42976	41328	3.16%	0.212%
53	11.233	1553	1555	1558	rVB	40803	33233	2.54%	0.171%
54	11.333	1568	1572	1575	rBV	1355014	1307780	100.00%	6.723%
55	11.357	1575	1576	1581	rVV	345968	274380	20.98%	1.411%
56	11.416	1583	1586	1590	rVB	96254	89412	6.84%	0.460%
57	11.580	1608	1614	1616	rBV4	29721	47259	3.61%	0.243%
58	11.610	1616	1619	1624	rVB2	57960	63972	4.89%	0.329%
59	11.669	1624	1629	1633	rBV	31138	36568	2.80%	0.188%
60	11.710	1633	1636	1640	rVB	128396	101105	7.73%	0.520%
61	11.757	1640	1644	1647	rBV2	15232	19813	1.52%	0.102%
62	11.839	1654	1658	1661	rBV	92610	104137	7.96%	0.535%
63	11.869	1661	1663	1667	rVB	108141	94873	7.25%	0.488%
64	11.916	1668	1671	1674	rVV	44235	43177	3.30%	0.222%
65	11.951	1674	1677	1679	rVV2	148112	151769	11.61%	0.780%
66	11.975	1679	1681	1685	rVB	65429	62926	4.81%	0.323%
67	12.016	1685	1688	1691	rBV	80800	69229	5.29%	0.356%
68	12.133	1705	1708	1712	rBV	51099	56915	4.35%	0.293%
69	12.180	1712	1716	1719	rVB4	24938	35516	2.72%	0.183%
70	12.263	1727	1730	1731	rBV3	20109	19050	1.46%	0.098%
71	12.310	1736	1738	1741	rVB	33883	30186	2.31%	0.155%

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72	12.339	1741	1743	1745	rVB	31501	21444	1.64%	0.110%
73	12.398	1748	1753	1755	rBV3	432662	470009	35.94%	2.416%
74	12.416	1755	1756	1760	rVV3	258559	231007	17.66%	1.188%
75	12.486	1763	1768	1769	rBV2	62990	76507	5.85%	0.393%
76	12.504	1769	1771	1776	rVB2	81949	74273	5.68%	0.382%
77	12.551	1776	1779	1783	rBV	636035	566859	43.35%	2.914%
78	12.592	1784	1786	1788	rVB3	24982	18313	1.40%	0.094%
79	12.616	1788	1790	1792	rBV2	24541	26627	2.04%	0.137%
80	12.704	1802	1805	1807	rBV	51871	49866	3.81%	0.256%
81	12.780	1812	1818	1824	rBV	598608	690165	52.77%	3.548%
82	12.833	1825	1827	1830	rVB	47428	40243	3.08%	0.207%
83	12.886	1834	1836	1837	rBV2	28952	24173	1.85%	0.124%
84	12.916	1837	1841	1844	rVV	749092	649134	49.64%	3.337%
85	12.998	1851	1855	1860	rBV2	59430	102369	7.83%	0.526%
86	13.110	1867	1874	1877	rBV	94335	181051	13.84%	0.931%
87	13.174	1883	1885	1888	rVB	52383	41381	3.16%	0.213%
88	13.216	1889	1892	1896	rBV2	83159	92517	7.07%	0.476%
89	13.310	1905	1908	1911	rBV	73671	63932	4.89%	0.329%
90	13.339	1911	1913	1916	rVB	79923	66773	5.11%	0.343%
91	13.733	1978	1980	1983	rBV2	45876	40533	3.10%	0.208%
92	13.804	1987	1992	1996	rBV3	75108	120386	9.21%	0.619%
93	13.986	2018	2023	2025	rBV2	884914	1070077	81.82%	5.501%
94	14.010	2025	2027	2030	rVB	247327	207855	15.89%	1.069%
95	14.092	2039	2041	2043	rBV	37674	33378	2.55%	0.172%
96	14.374	2087	2089	2092	rVB	80478	66962	5.12%	0.344%
97	15.027	2197	2200	2203	rBV	239018	330576	25.28%	1.699%
98	15.327	2248	2251	2255	rBV	164800	192990	14.76%	0.992%
99	15.392	2259	2262	2266	rBV	229532	261843	20.02%	1.346%
100	15.457	2269	2273	2276	rVB	631678	724510	55.40%	3.725%

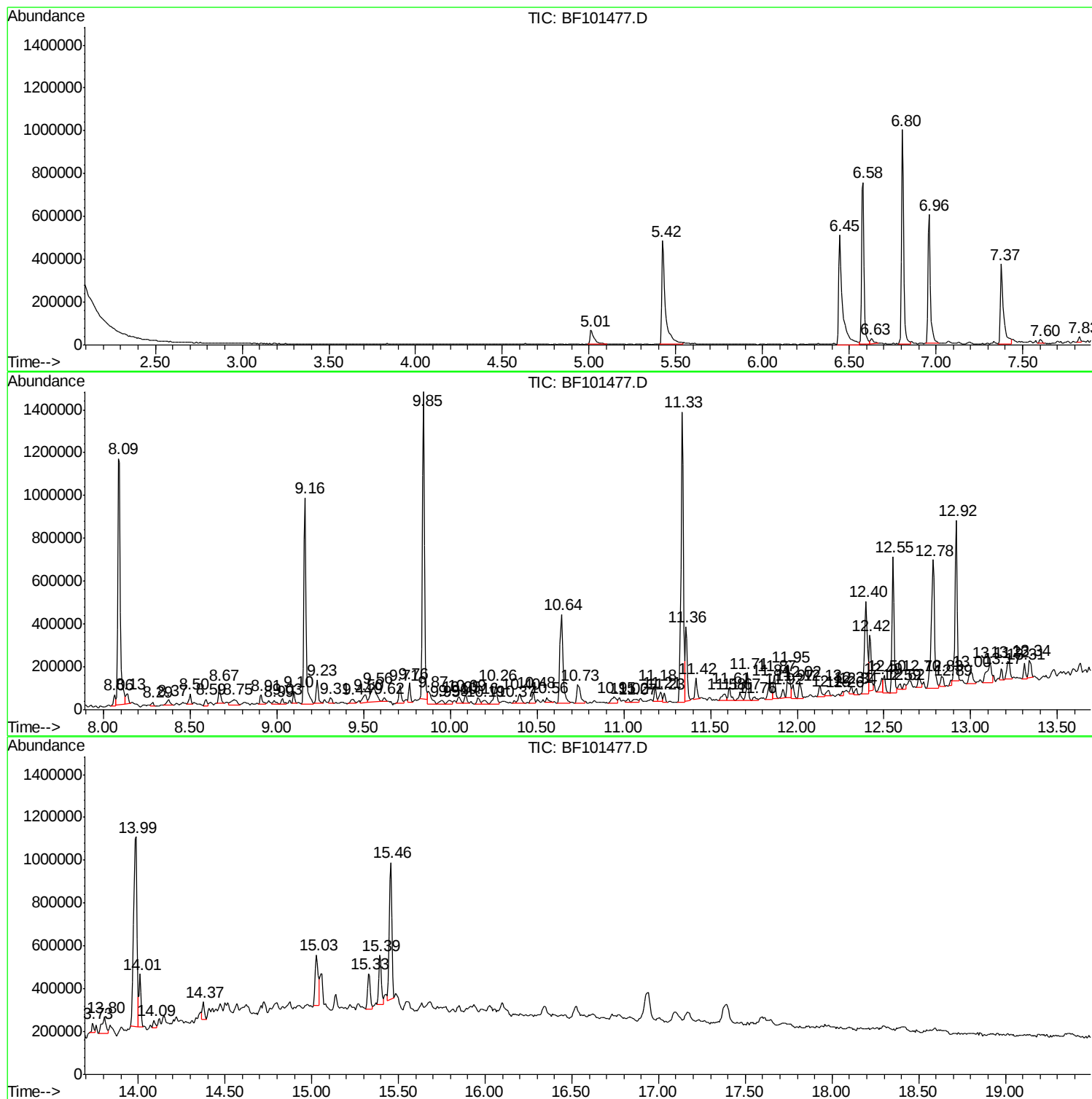
Sum of corrected areas: 19451913

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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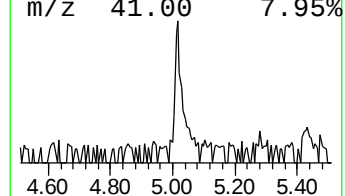
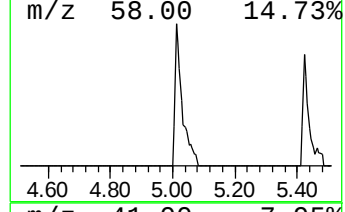
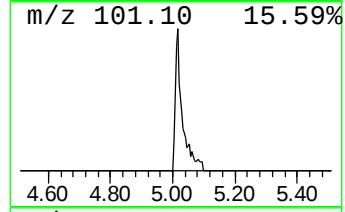
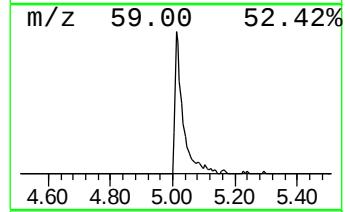
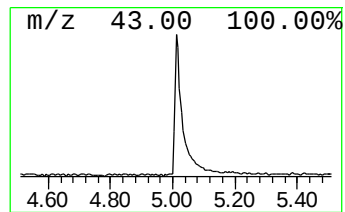
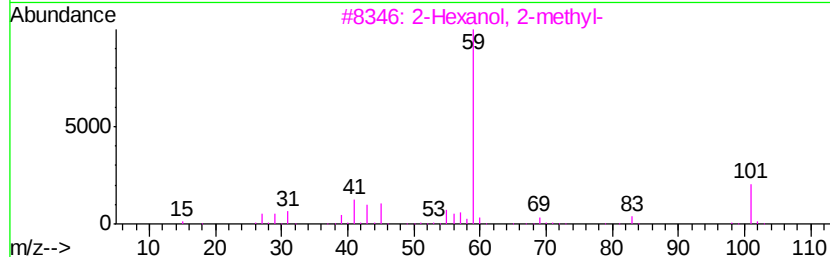
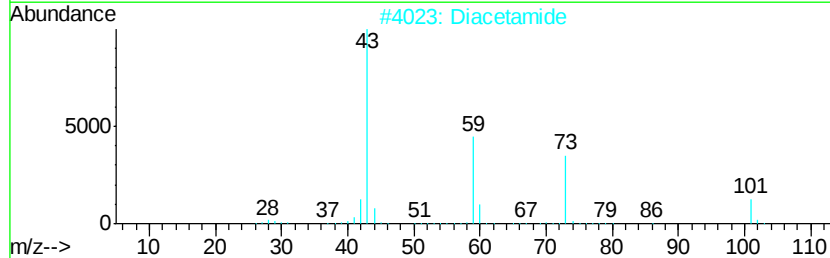
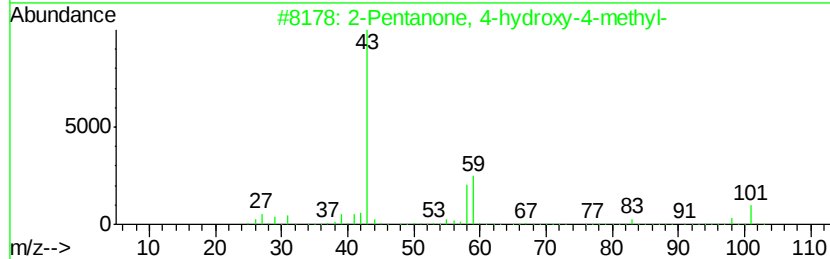
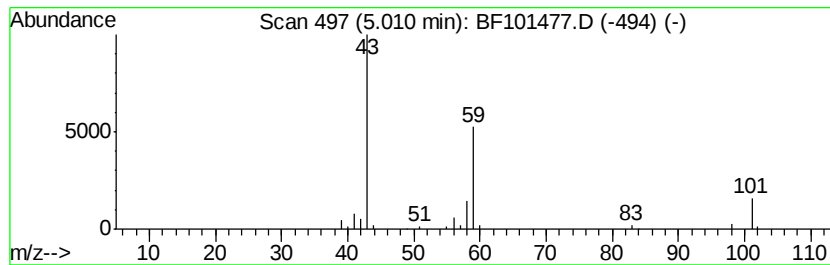
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	2.33 ng	108870	1,4-Dichlorobenzene-d4	6.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Diacetamide	101	C4H7NO2	000625-77-4	36	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9	



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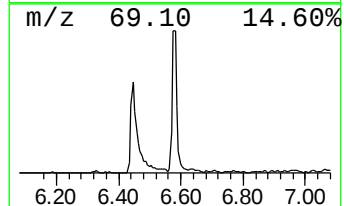
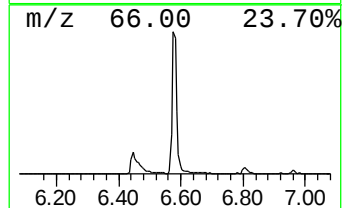
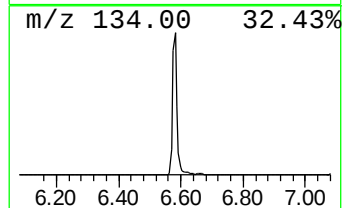
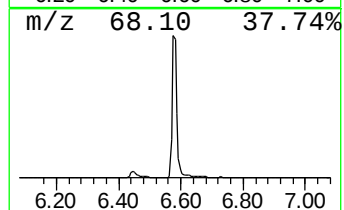
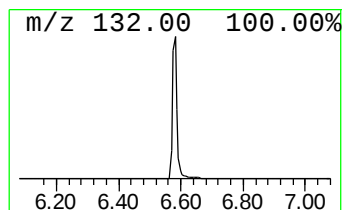
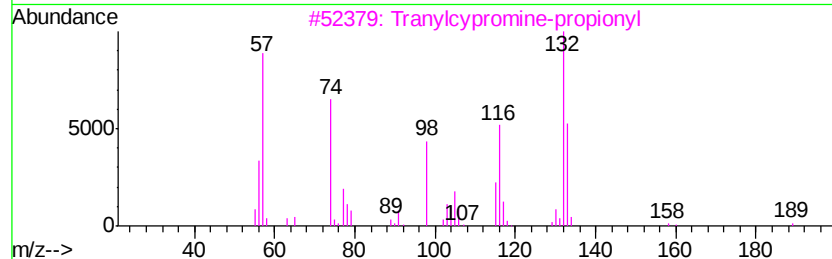
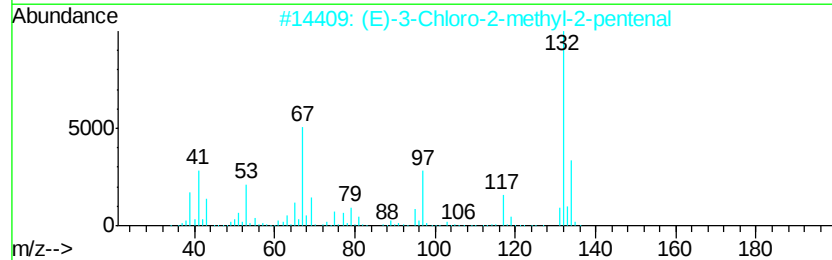
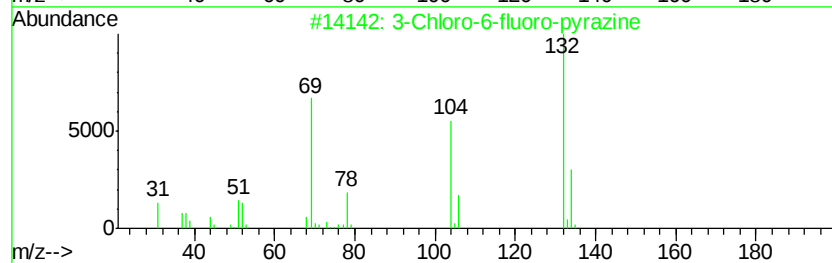
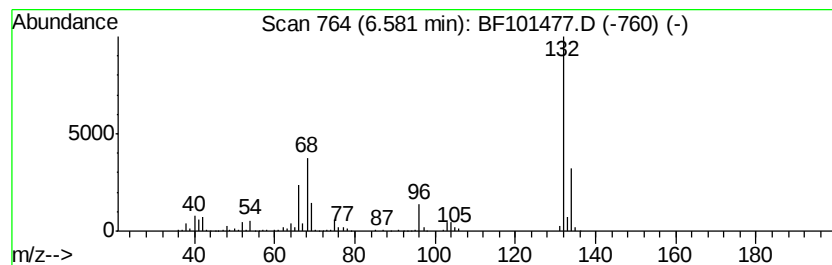
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	16.90 ng	791283	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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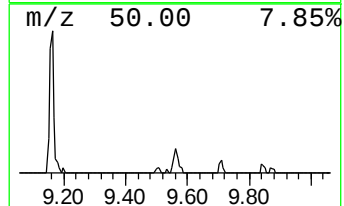
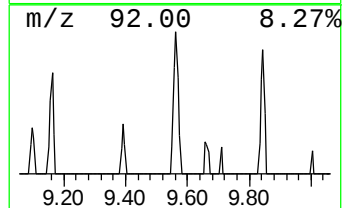
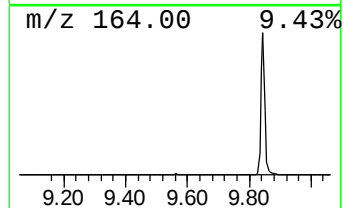
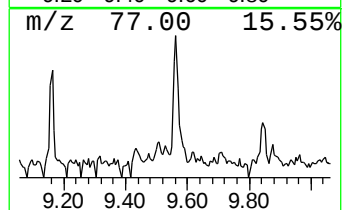
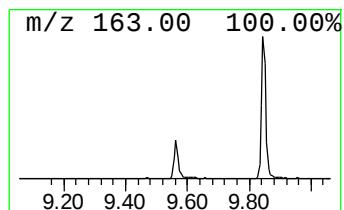
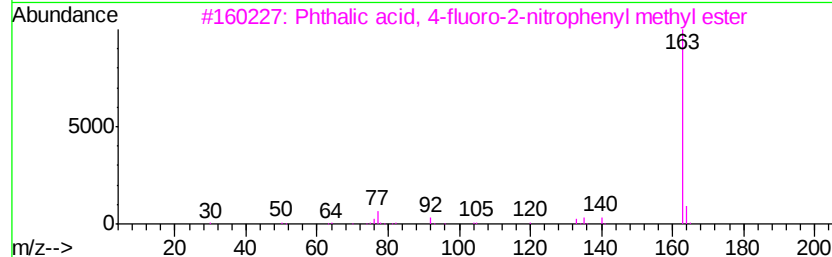
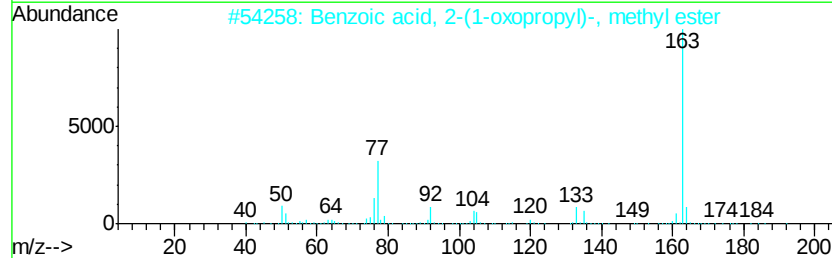
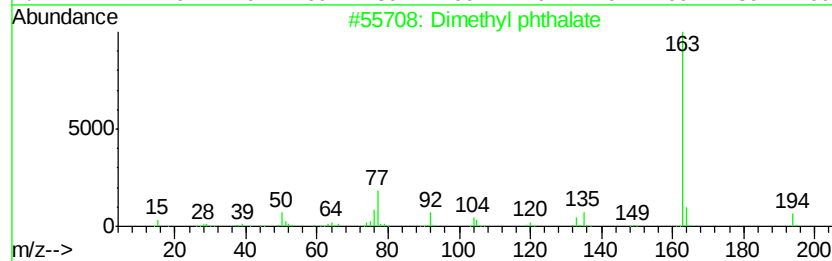
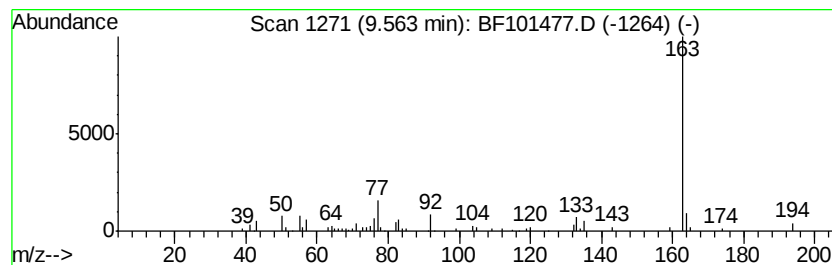
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Dimethyl phthalate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	2.31 ng	150270	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl phthalate	194	C10H10O4	000131-11-3	50
2		Benzoic acid, 2-(1-oxopropyl)-, ...	192	C11H12O3	032025-37-9	50
3		Phthalic acid, 4-fluoro-2-nitrop...	319	C15H10FN06	1000315-63-4	50
4		2-Methoxyethyl methyl phthalate	238	C12H14O5	1000373-89-4	47
5		Methyl pentyl phthalate	250	C14H18O4	1000373-89-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
Data File : BF101477.D
Acq On : 18 Dec 2017 13:42
Operator : SJ/JU
Sample : I6679-07 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HR-05-121317-C

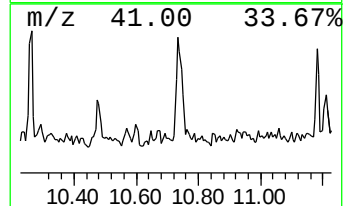
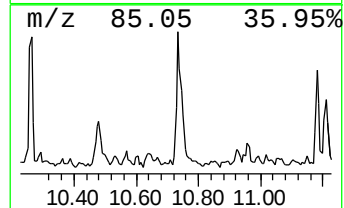
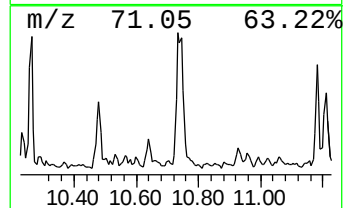
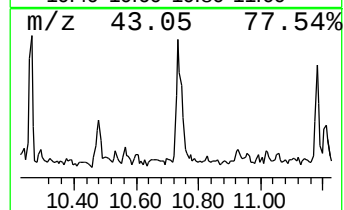
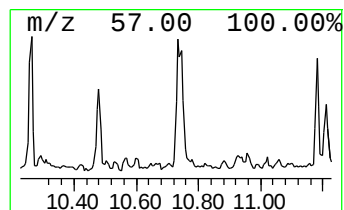
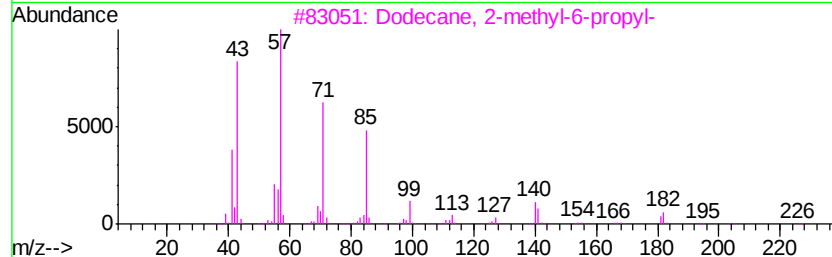
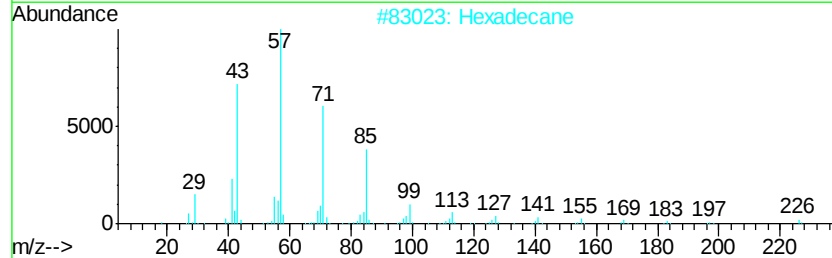
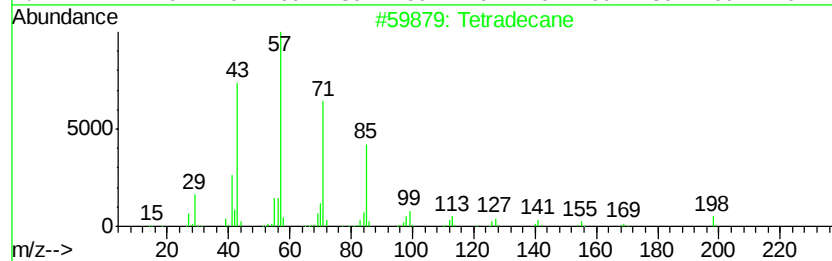
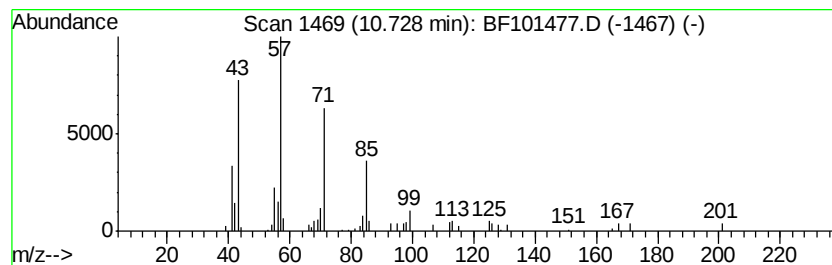
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tetradecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	2.02 ng	132094	Phenanthrene-d10	11.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	90
2		Hexadecane	226	C16H34	000544-76-3	90
3		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4		Octadecane	254	C18H38	000593-45-3	86
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
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Acq On : 18 Dec 2017 13:42
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Sample : I6679-07 5X
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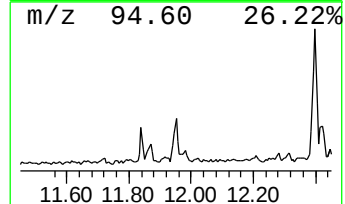
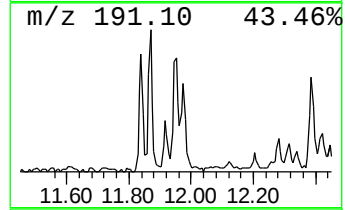
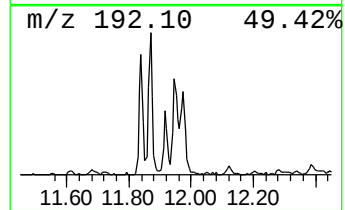
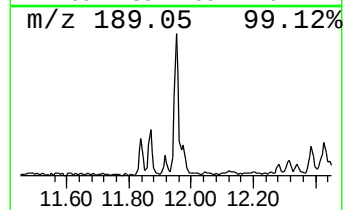
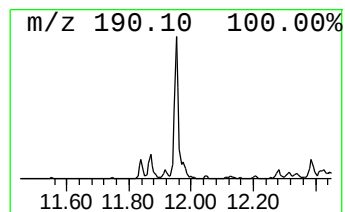
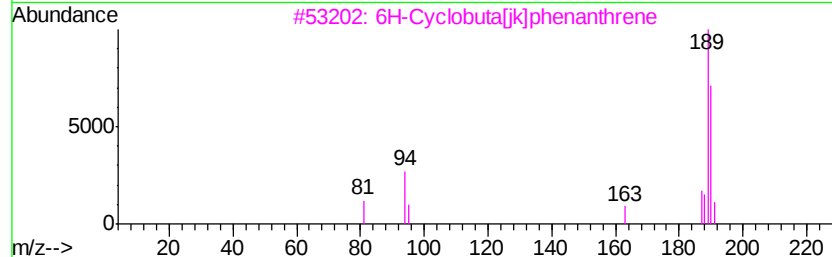
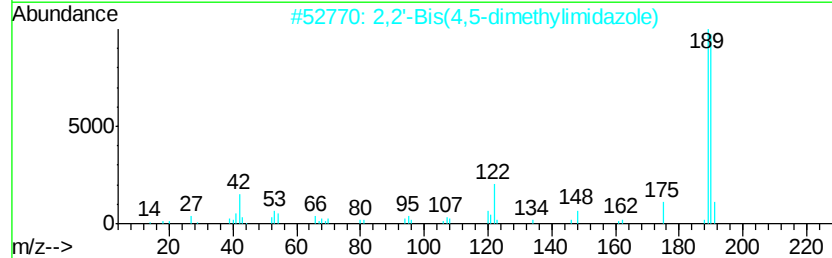
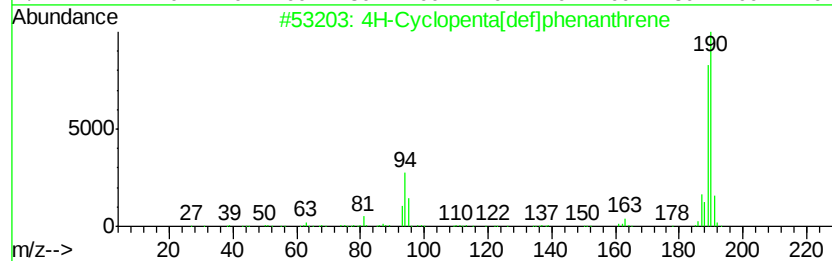
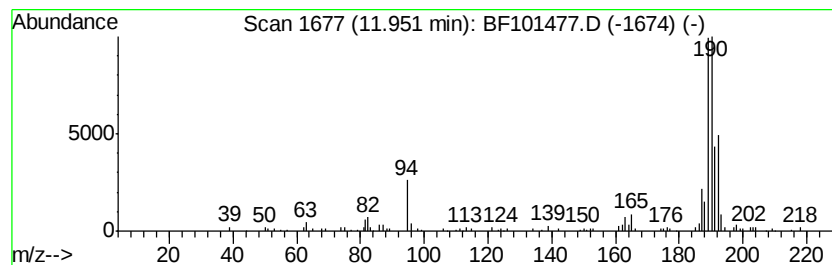
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	2.32 ng	151769	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
4	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	33
5	1,1'-Biphenyl, 4,4'-difluoro-	190	C12H8F2	000398-23-2	23



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
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Sample : I6679-07 5X
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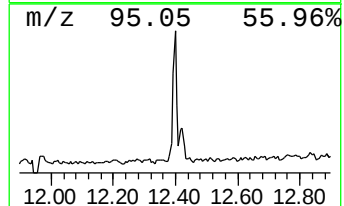
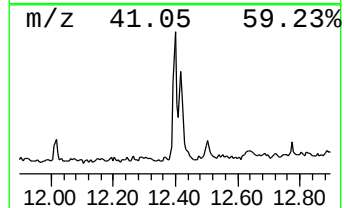
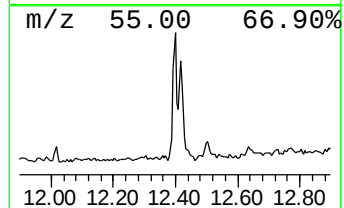
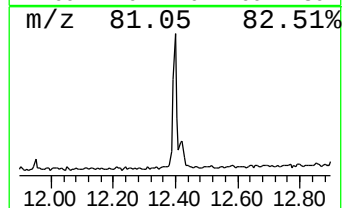
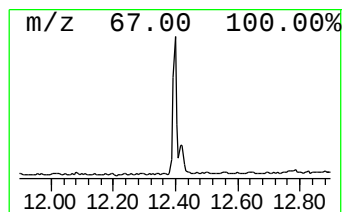
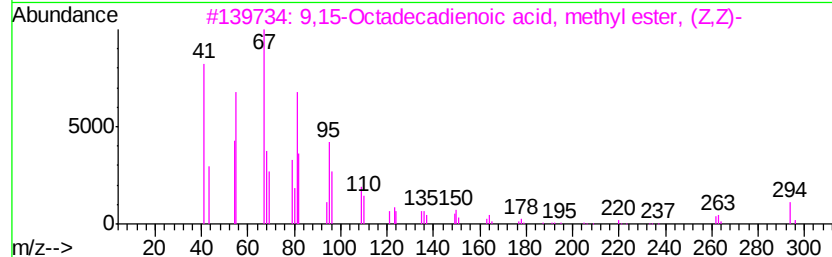
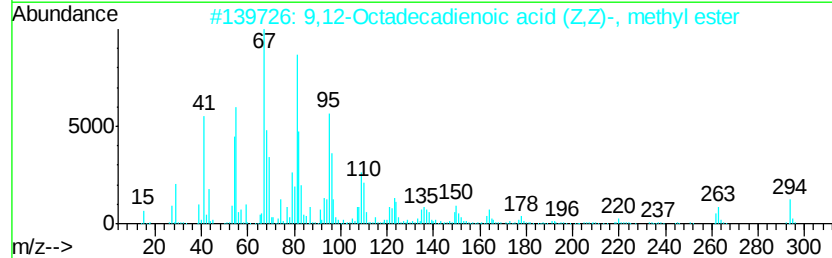
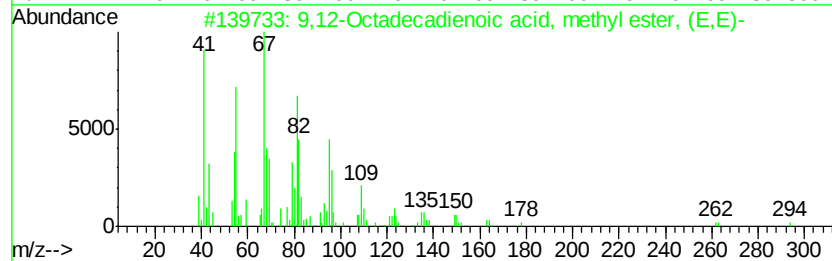
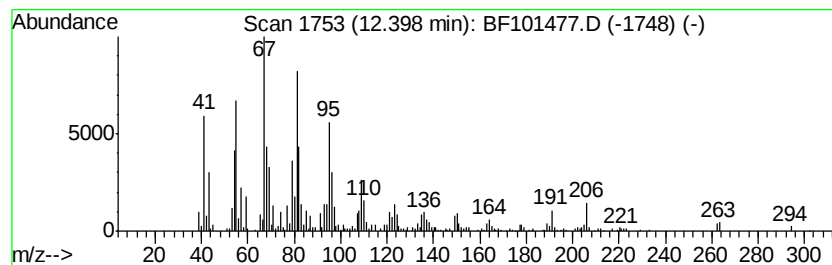
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	7.19 ng	470009	Phenanthrene-d10	11.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3			9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
5			11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
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Operator : SJ/JU
Sample : I6679-07 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

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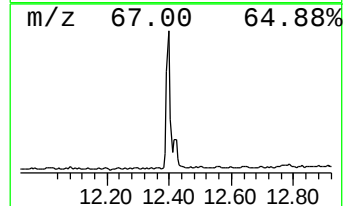
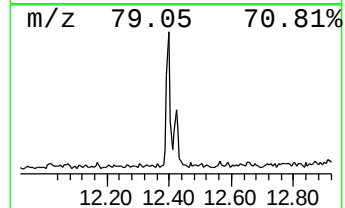
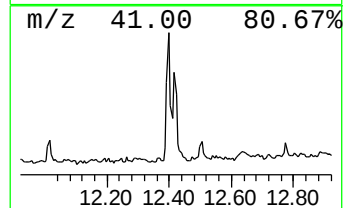
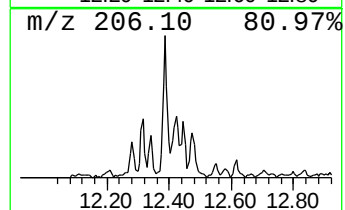
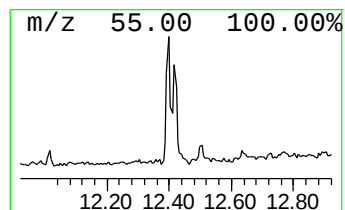
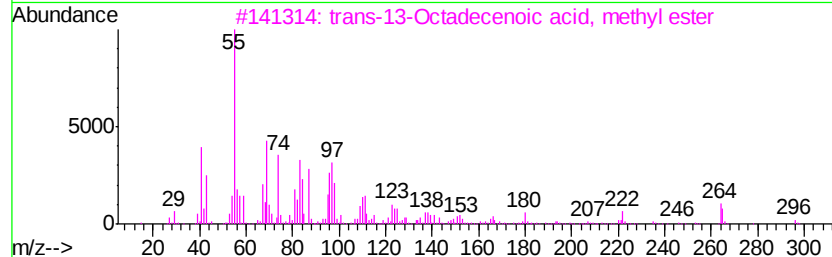
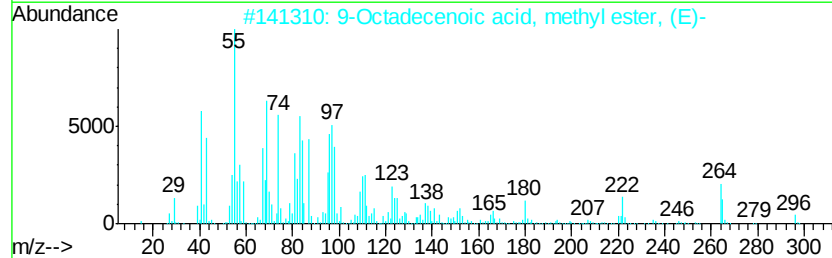
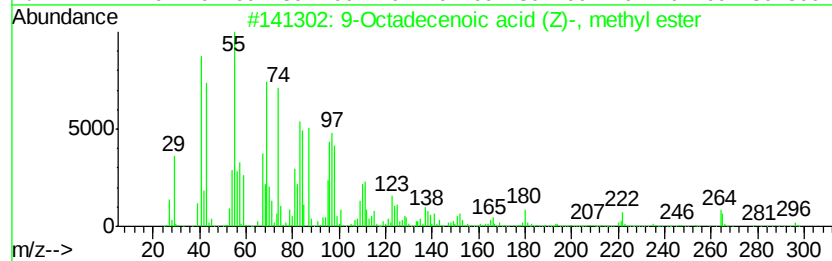
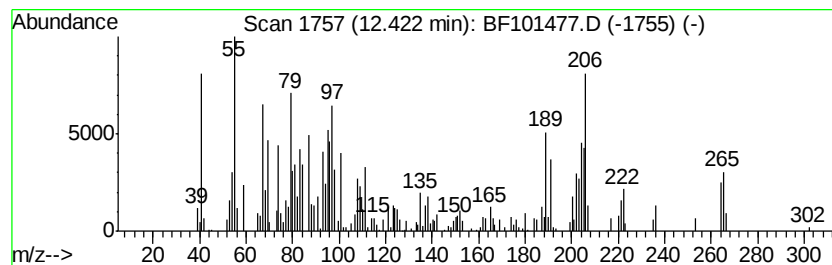
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenoic acid (Z)-, m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	3.53 ng	231007	Phenanthrene-d10	11.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	93
2		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	76
3		trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	74
4		cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	72
5		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
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Misc :
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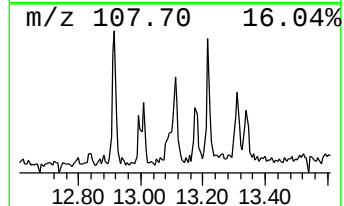
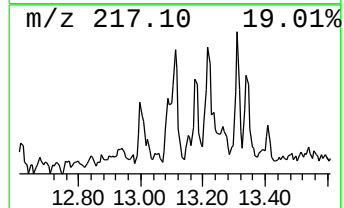
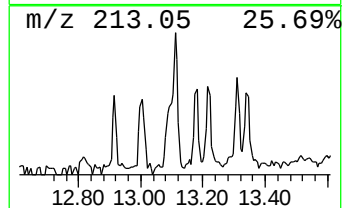
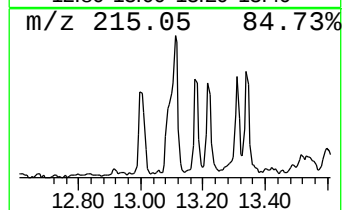
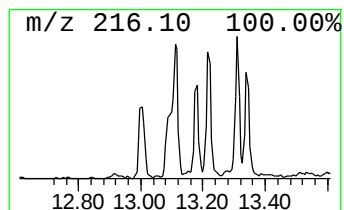
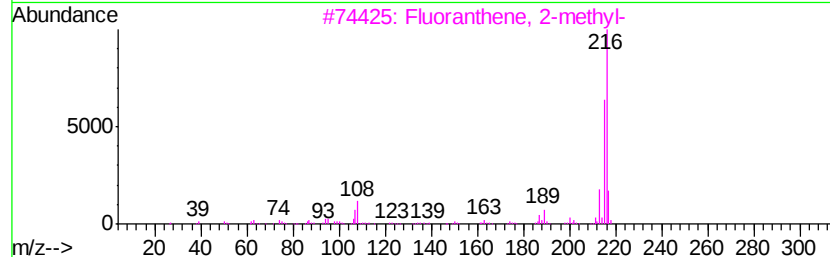
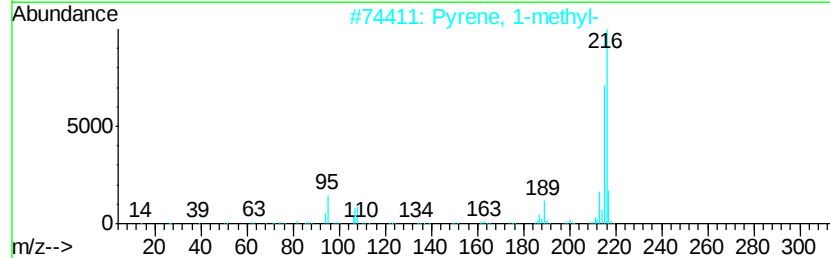
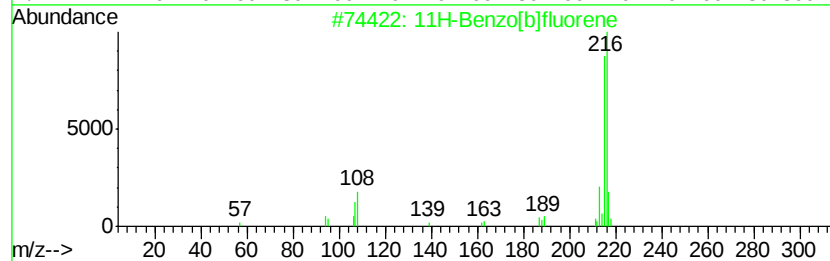
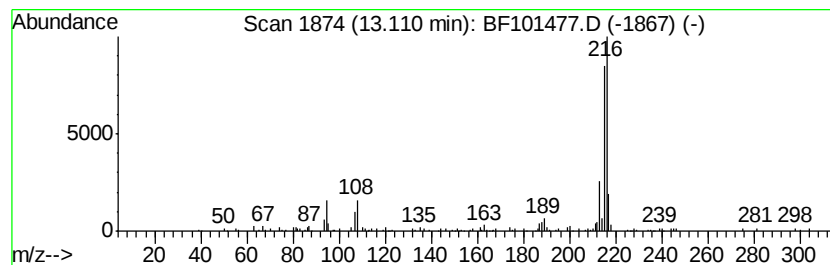
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	3.38 ng	181051	Chrysene-d12	13.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



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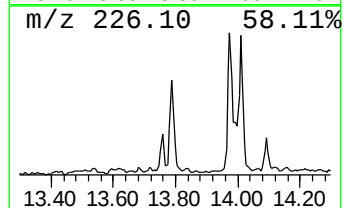
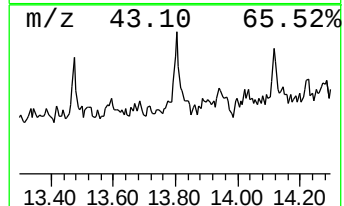
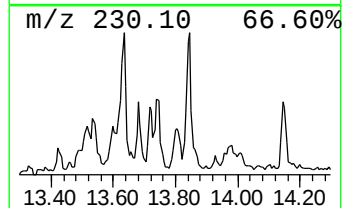
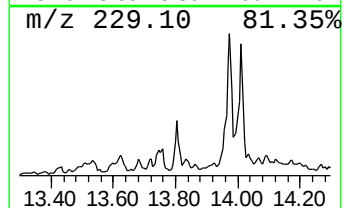
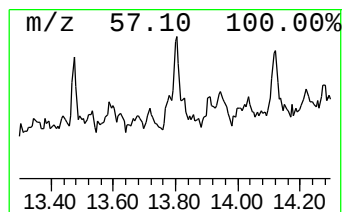
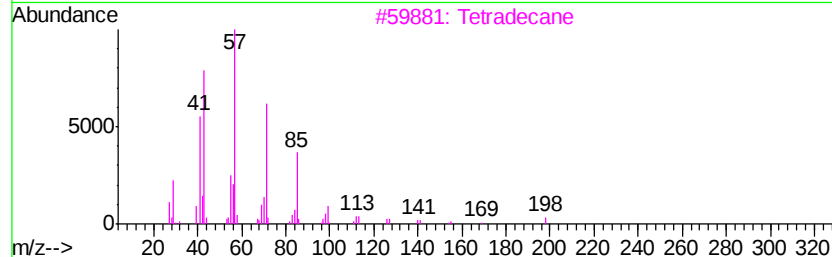
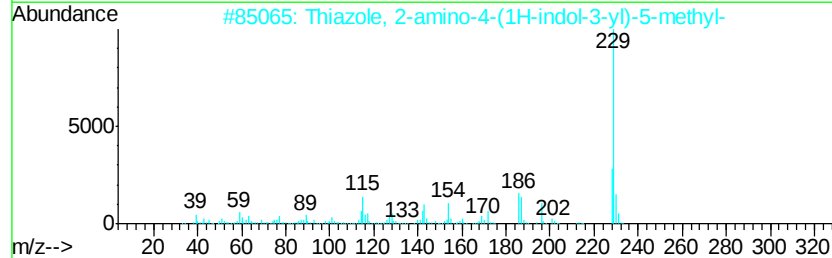
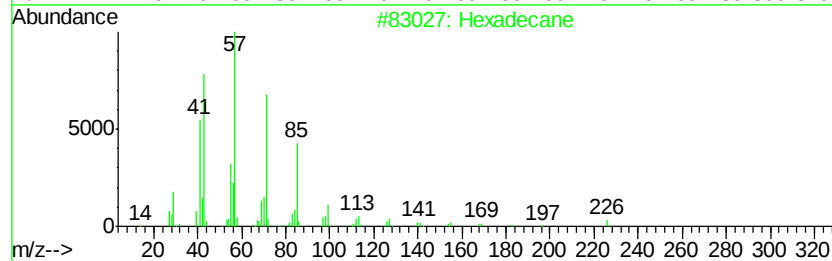
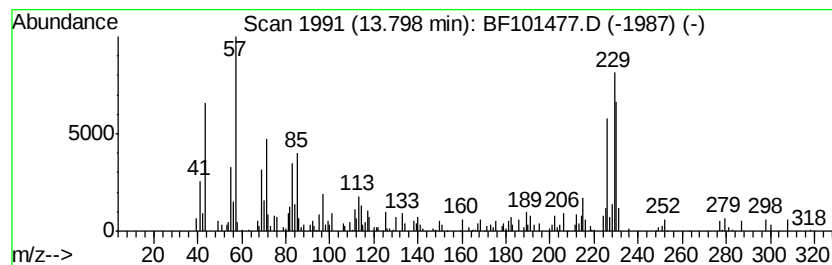
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	2.25 ng	120386	Chrysene-d12	13.99
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane		226 C16H34	000544-76-3 44
2	Thiazole, 2-amino-4-(1H-indol-3-yl)-5-methyl-		229 C12H11N3S	1000304-18-3 38
3	Tetradecane		198 C14H30	000629-59-4 25
4	Tridecane		184 C13H28	000629-50-5 25
5	Sulfurous acid, octadecyl 2-prop...		376 C21H44O3S	1000309-12-7 20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
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Sample : I6679-07 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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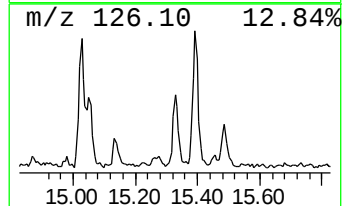
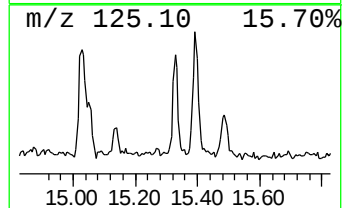
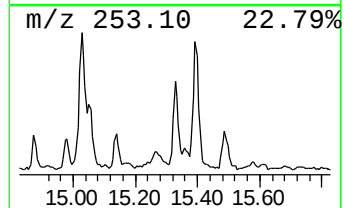
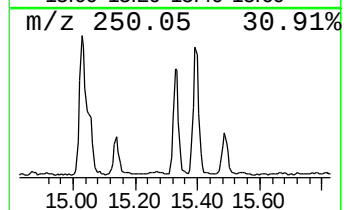
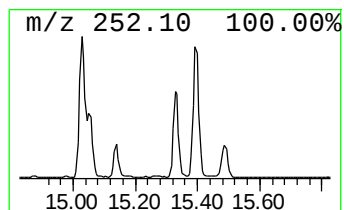
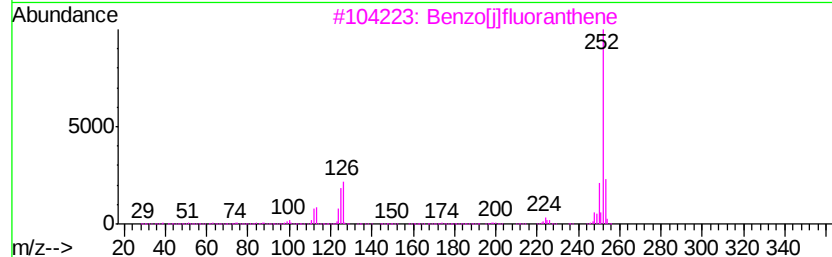
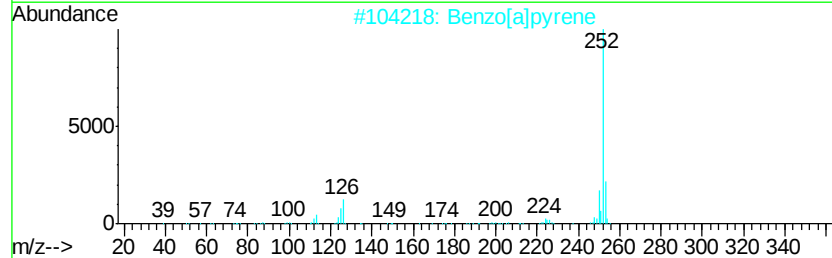
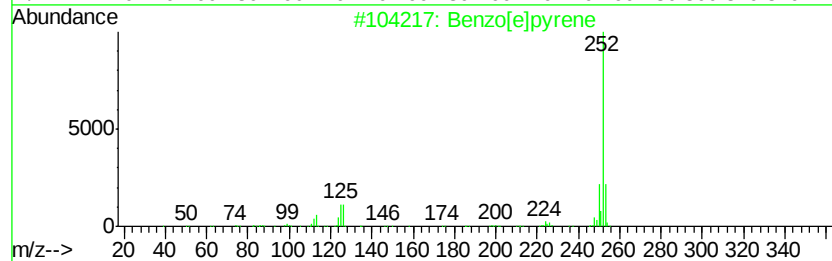
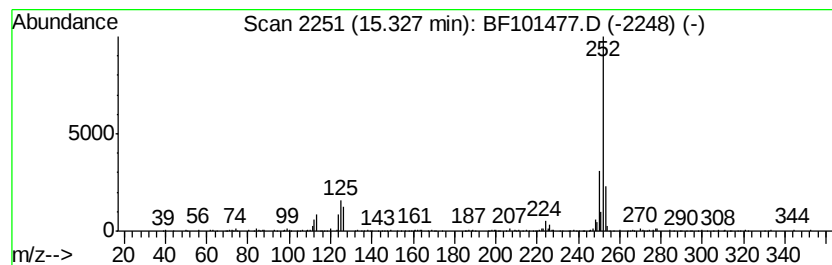
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	5.33 ng	192990	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2		Benzo[a]pyrene	252	C20H12	000050-32-8	96
3		Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
Data File : BF101477.D
Acq On : 18 Dec 2017 13:42
Operator : SJ/JU
Sample : I6679-07 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HR-05-121317-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	2.3	ng	108870	1	6.80	936162	20.0
unknown6.58	6.58	16.9	ng	791283	1	6.80	936162	20.0
Dimethyl phthalate	9.56	2.3	ng	150270	3	9.85	1302310	20.0
Tetradecane	10.73	2.0	ng	132094	4	11.33	1307780	20.0
4H-Cyclopenta[def...	11.95	2.3	ng	151769	4	11.33	1307780	20.0
9,12-Octadecadien...	12.40	7.2	ng	470009	4	11.33	1307780	20.0
9-Octadecenoic ac...	12.42	3.5	ng	231007	4	11.33	1307780	20.0
11H-Benzo[b]fluorene	13.11	3.4	ng	181051	5	13.99	1070080	20.0
Hexadecane	13.80	2.3	ng	120386	5	13.99	1070080	20.0
Benzo[e]pyrene	15.33	5.3	ng	192990	6	15.46	724510	20.0