

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
 Data File : BF096356.D
 Acq On : 1 Jul 2017 00:32
 Operator : SJ/MA
 Sample : I3930-03
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B2-4-6

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-BF062717.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	2.846	123	129	137	rVB3	21477	42225	1.06%	0.094%
2	3.787	282	289	299	rVB	1718035	2542136	63.89%	5.662%
3	4.393	387	392	398	rVB4	38600	62222	1.56%	0.139%
4	4.963	484	489	496	rBV	495564	694194	17.45%	1.546%
5	5.345	548	554	555	rBV	46823	46663	1.17%	0.104%
6	5.381	555	560	563	rVB	3677753	3933744	98.86%	8.761%
7	6.210	696	701	705	rBV4	23037	42772	1.07%	0.095%
8	6.298	713	716	721	rVB3	46649	58109	1.46%	0.129%
9	6.357	721	726	727	rBV	45367	48126	1.21%	0.107%
10	6.392	727	732	736	rVB	3404813	3979136	100.00%	8.862%
11	6.528	750	755	759	rBV	3590261	3735423	93.88%	8.320%
12	6.592	762	766	767	rBV	46938	44125	1.11%	0.098%
13	6.757	791	794	797	rBV	1051170	860858	21.63%	1.917%
14	6.822	802	805	813	rBV2	50875	64283	1.62%	0.143%
15	6.916	817	821	824	rBV	2990148	2715776	68.25%	6.049%
16	7.163	857	863	865	rBV3	42487	45360	1.14%	0.101%
17	7.316	881	889	892	rBV	2354583	2447883	61.52%	5.452%
18	7.369	895	898	901	rVV2	67069	59379	1.49%	0.132%
19	7.410	901	905	908	rVB3	59989	71770	1.80%	0.160%
20	8.039	1008	1012	1015	rBV	1350395	1207659	30.35%	2.690%
21	8.063	1015	1016	1018	rVB	161350	73586	1.85%	0.164%
22	8.381	1066	1070	1074	rBV4	32889	45794	1.15%	0.102%
23	8.428	1074	1078	1081	rVV3	73798	82789	2.08%	0.184%
24	8.481	1083	1087	1090	rVB	68602	65859	1.66%	0.147%
25	8.645	1112	1115	1117	rBV	66892	50862	1.28%	0.113%
26	8.751	1127	1133	1136	rVB	161244	138742	3.49%	0.309%
27	8.851	1147	1150	1154	rVB	122150	94020	2.36%	0.209%
28	9.116	1190	1195	1199	rVB	3077012	2857492	71.81%	6.364%
29	9.210	1208	1211	1215	rBV2	110169	103200	2.59%	0.230%
30	9.375	1236	1239	1244	rVB2	46060	52915	1.33%	0.118%
31	9.451	1250	1252	1255	rBV	79157	73433	1.85%	0.164%
32	9.480	1255	1257	1259	rVV	69359	57596	1.45%	0.128%
33	9.510	1259	1262	1264	rVV	572939	463704	11.65%	1.033%
34	9.533	1264	1266	1268	rVB	56207	40860	1.03%	0.091%

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35	9.651	1282	1286	1290	rBV	214513	175893	4.42%	0.392%
36	9.739	1298	1301	1305	rVB	134677	107318	2.70%	0.239%
37	9.792	1305	1310	1313	rBV	1075611	965238	24.26%	2.150%
38	9.822	1313	1315	1319	rVB3	44423	43330	1.09%	0.097%
39	9.998	1342	1345	1348	rVB	134962	119067	2.99%	0.265%
40	10.204	1376	1380	1383	rBV3	52707	75961	1.91%	0.169%
41	10.239	1383	1386	1389	rVB	140003	122490	3.08%	0.273%
42	10.333	1399	1402	1405	rBV2	70297	57399	1.44%	0.128%
43	10.457	1419	1423	1426	rBV3	65810	78650	1.98%	0.175%
44	10.492	1426	1429	1435	rVB3	42130	68665	1.73%	0.153%
45	10.575	1439	1443	1447	rVB	1667682	1628674	40.93%	3.627%
46	10.710	1461	1466	1473	rBV3	200337	331202	8.32%	0.738%
47	11.075	1525	1528	1533	rVB	74467	71010	1.78%	0.158%
48	11.127	1533	1537	1539	rVV4	30443	43544	1.09%	0.097%
49	11.157	1539	1542	1545	rVV2	116612	134723	3.39%	0.300%
50	11.186	1545	1547	1551	rVB2	41859	41518	1.04%	0.092%
51	11.275	1558	1562	1564	rBV	859370	803660	20.20%	1.790%
52	11.298	1564	1566	1569	rVV	443201	350356	8.80%	0.780%
53	11.345	1569	1574	1577	rVV	272055	237391	5.97%	0.529%
54	11.380	1577	1580	1585	rVB6	35555	43835	1.10%	0.098%
55	11.498	1597	1600	1603	rVB2	71186	61523	1.55%	0.137%
56	11.580	1611	1614	1616	rBV2	60228	51667	1.30%	0.115%
57	11.774	1644	1647	1649	rBV	85389	71462	1.80%	0.159%
58	11.816	1649	1654	1656	rVV2	456682	477509	12.00%	1.064%
59	11.863	1656	1662	1664	rVV	488873	789795	19.85%	1.759%
60	11.880	1664	1665	1668	rVB	81252	64785	1.63%	0.144%
61	11.986	1680	1683	1689	rVB4	68139	71635	1.80%	0.160%
62	12.069	1694	1697	1698	rVV2	80407	68082	1.71%	0.152%
63	12.086	1698	1700	1704	rVB2	62512	54910	1.38%	0.122%
64	12.216	1718	1722	1725	rBV5	44577	61665	1.55%	0.137%
65	12.321	1736	1740	1743	rBV2	109041	121779	3.06%	0.271%
66	12.374	1747	1749	1752	rVV2	91573	87479	2.20%	0.195%
67	12.404	1752	1754	1759	rVB2	81828	83723	2.10%	0.186%
68	12.480	1763	1767	1771	rBV	899422	788604	19.82%	1.756%
69	12.516	1771	1773	1780	rBV5	54863	96370	2.42%	0.215%
70	12.592	1784	1786	1791	rVV2	125296	113283	2.85%	0.252%
71	12.680	1798	1801	1802	rBV	106136	110560	2.78%	0.246%

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72	12.710	1802	1806	1809	rVB	876650	808523	20.32%	1.801%
73	12.827	1824	1826	1827	rBV	69648	48016	1.21%	0.107%
74	12.857	1827	1831	1834	rVV	2158202	2025352	50.90%	4.511%
75	12.886	1834	1836	1839	rVB	82921	69061	1.74%	0.154%
76	12.927	1840	1843	1846	rBV2	78343	90269	2.27%	0.201%
77	12.992	1852	1854	1856	rBV3	43617	48557	1.22%	0.108%
78	13.021	1856	1859	1860	rBV3	46815	59338	1.49%	0.132%
79	13.104	1870	1873	1875	rBV	101164	90419	2.27%	0.201%
80	13.139	1876	1879	1882	rBV	148954	139506	3.51%	0.311%
81	13.233	1892	1895	1897	rBV	80050	70710	1.78%	0.157%
82	13.263	1898	1900	1904	rVB3	69320	65869	1.66%	0.147%
83	13.339	1910	1913	1920	rVB	310834	334895	8.42%	0.746%
84	13.410	1922	1925	1928	rBV2	161359	176722	4.44%	0.394%
85	13.439	1928	1930	1933	rVB	64144	61915	1.56%	0.138%
86	13.539	1944	1947	1952	rBV	121847	147255	3.70%	0.328%
87	13.651	1962	1966	1969	rBV2	123686	175672	4.41%	0.391%
88	13.680	1969	1971	1973	rVV	89346	87026	2.19%	0.194%
89	13.704	1973	1975	1979	rVV2	108578	133924	3.37%	0.298%
90	13.757	1980	1984	1990	rVB2	311683	368310	9.26%	0.820%
91	13.904	2004	2009	2012	rBV2	1017019	1474323	37.05%	3.284%
92	13.927	2012	2013	2019	rVB	518052	410199	10.31%	0.914%
93	14.086	2038	2040	2042	rVB3	82905	61299	1.54%	0.137%
94	14.921	2178	2182	2185	rBV	550488	807696	20.30%	1.799%
95	15.210	2227	2231	2236	rVB	339556	442810	11.13%	0.986%
96	15.262	2237	2240	2244	rBV	317851	367577	9.24%	0.819%
97	15.327	2247	2251	2254	rBV	413337	480375	12.07%	1.070%
98	16.704	2482	2485	2492	rVB2	154114	275814	6.93%	0.614%

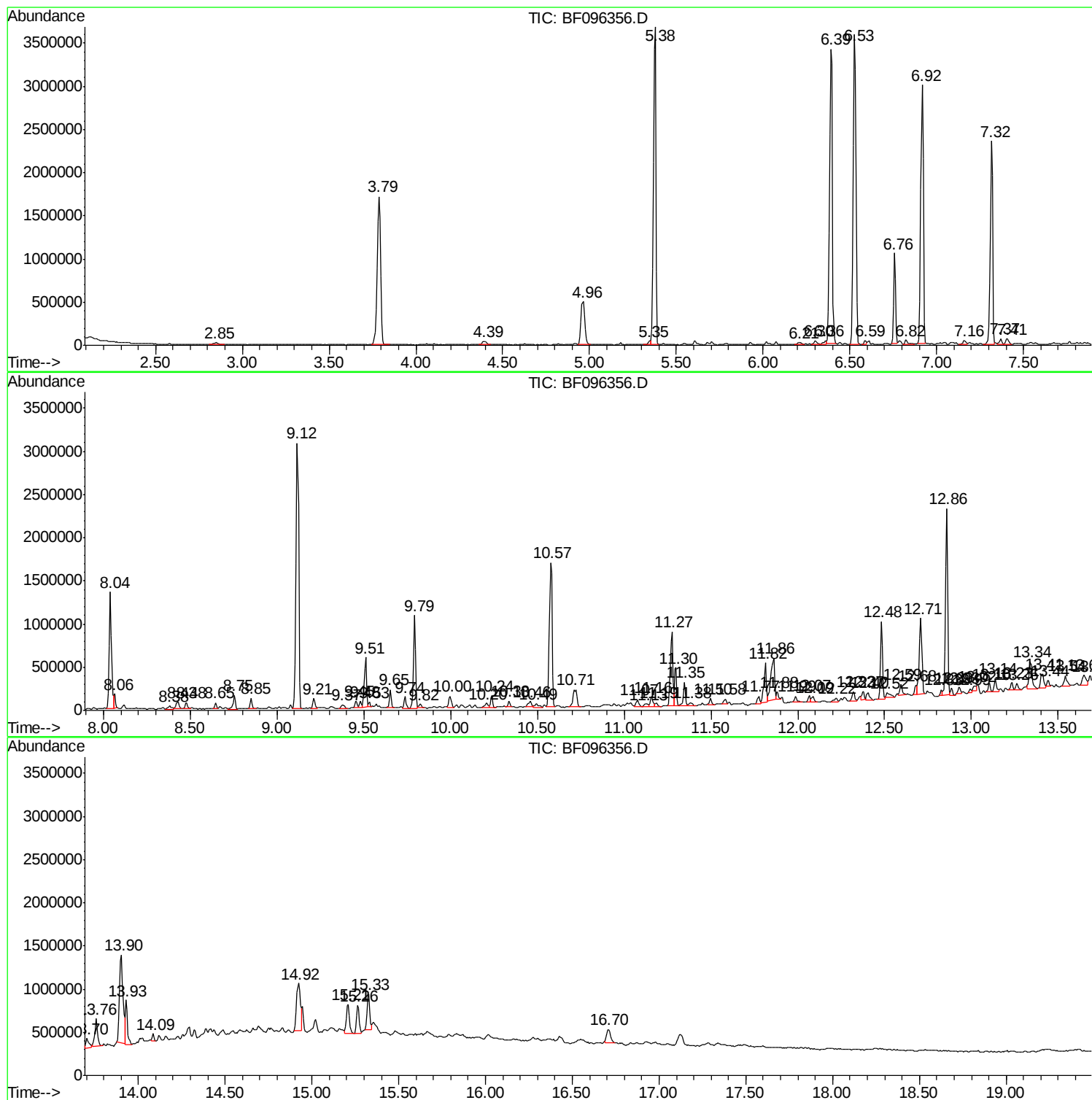
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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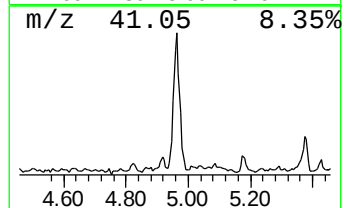
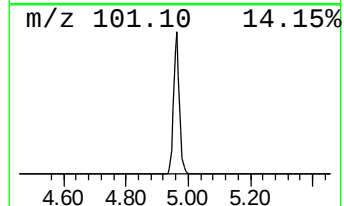
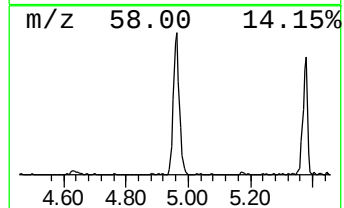
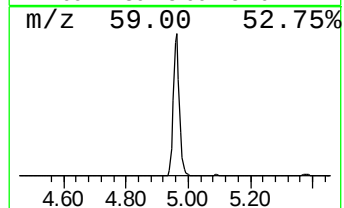
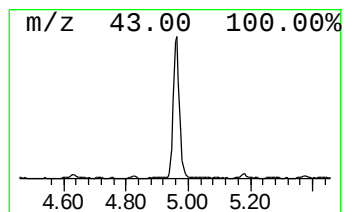
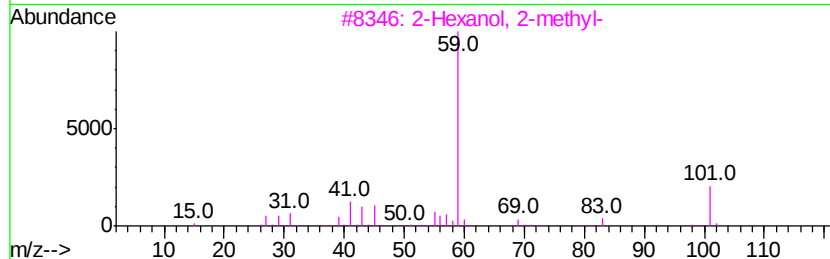
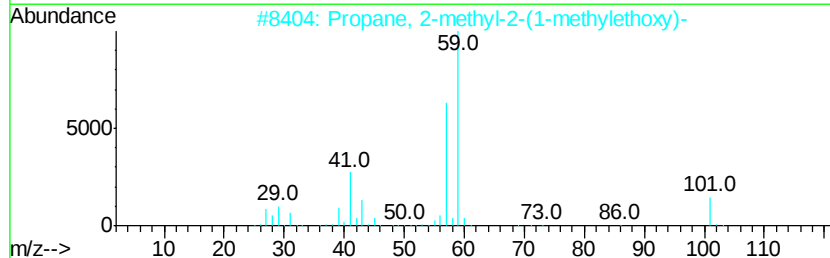
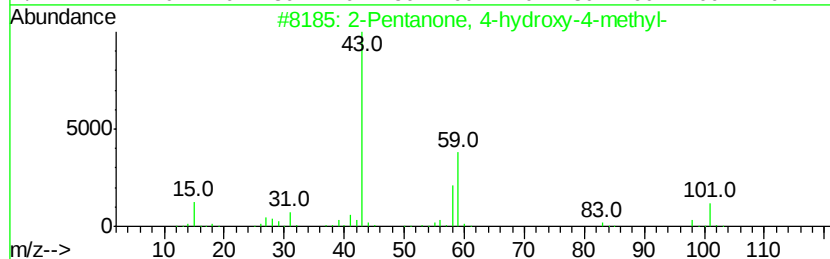
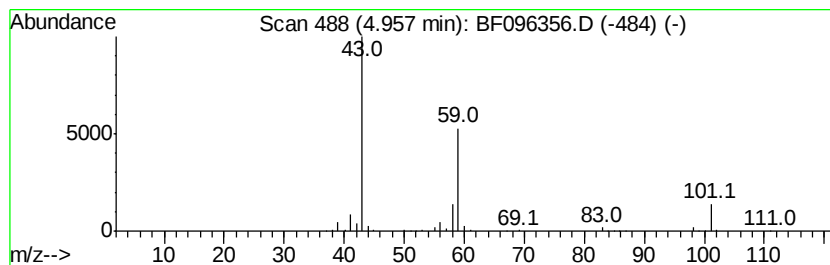
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	16.13 ng	694194	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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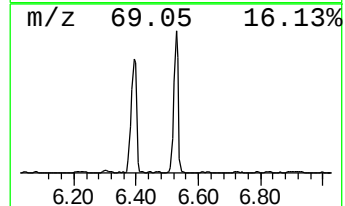
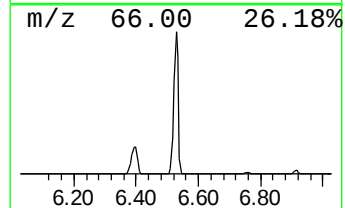
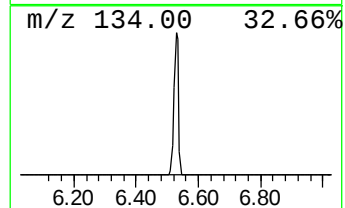
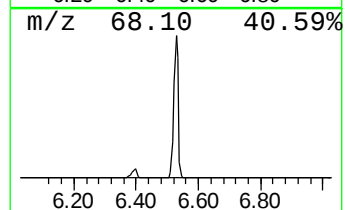
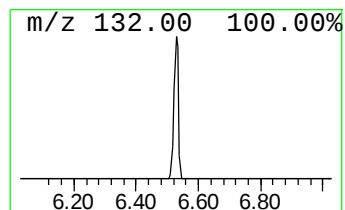
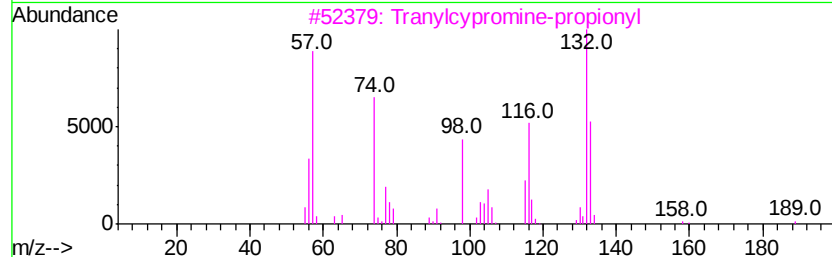
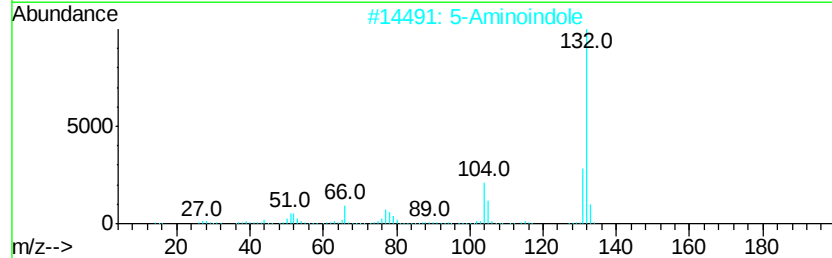
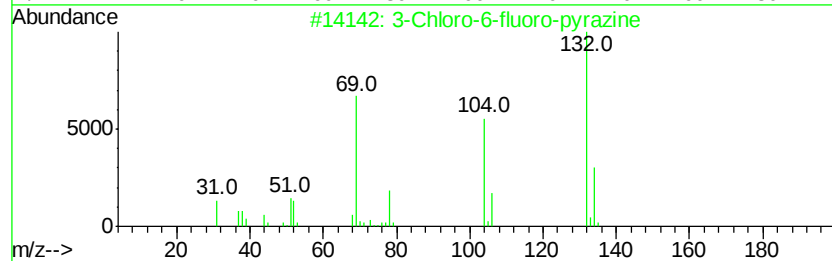
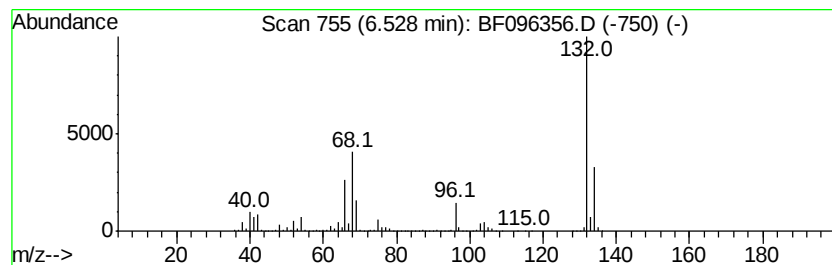
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	86.78 ng	3735420	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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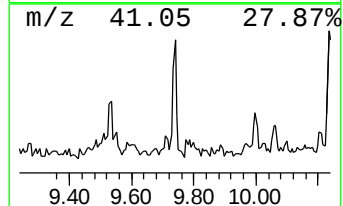
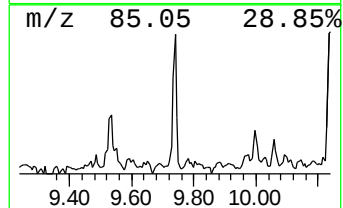
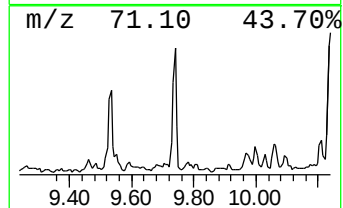
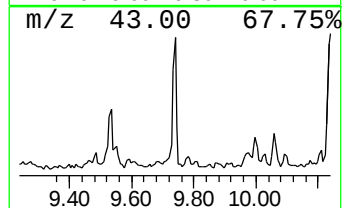
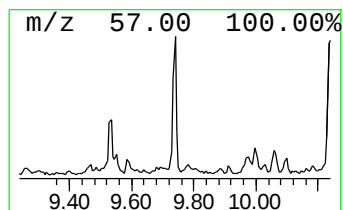
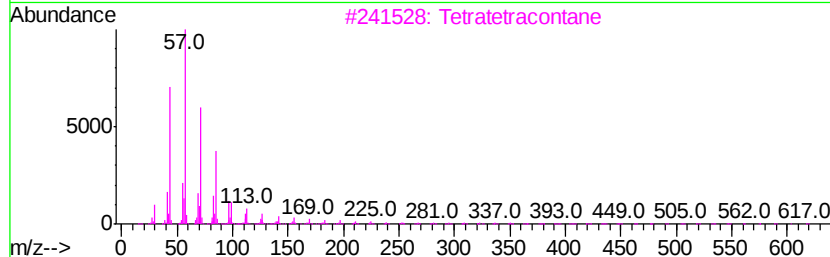
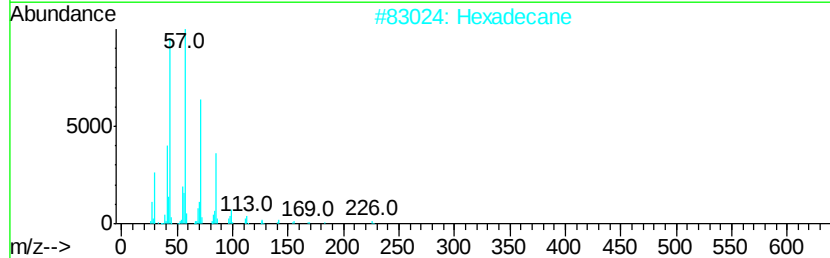
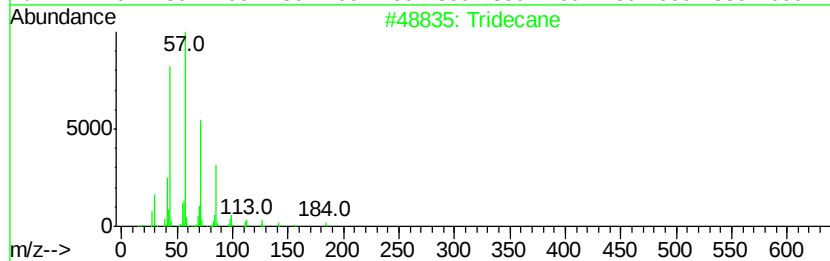
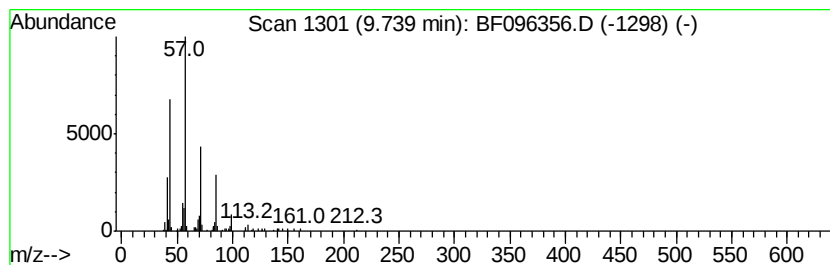
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Peak Number 6 Tridecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	2.22 ng	107318	Acenaphthene-d10	9.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	86
2	Hexadecane		226	C16H34	000544-76-3	80
3	Tetratetracontane		619	C44H90	007098-22-8	78
4	Tetradecane		198	C14H30	000629-59-4	78
5	Tridecane, 4,8-dimethyl-		212	C15H32	055030-62-1	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
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Acq On : 1 Jul 2017 00:32
Operator : SJ/MA
Sample : I3930-03
Misc :
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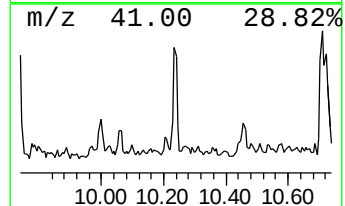
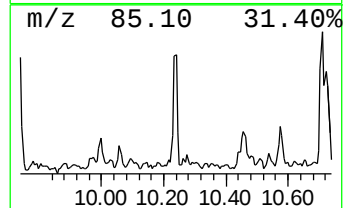
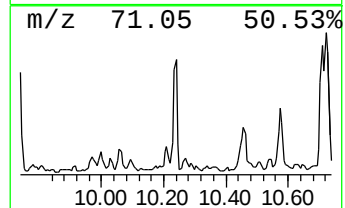
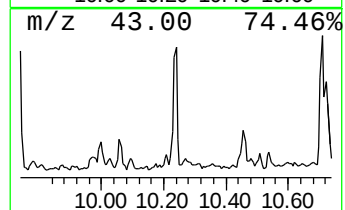
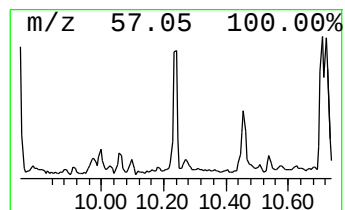
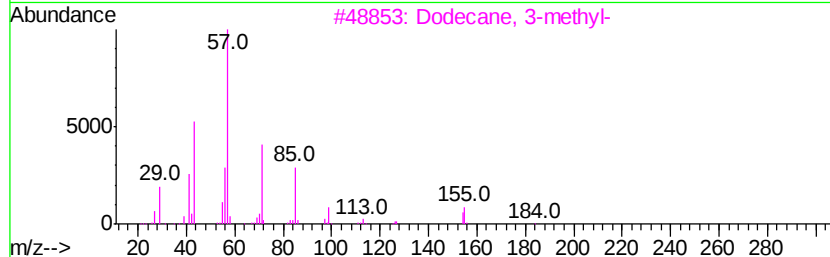
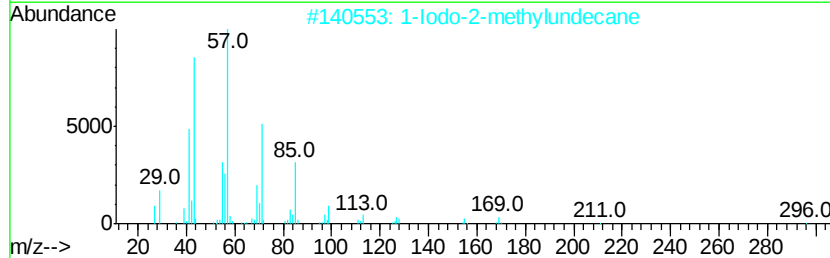
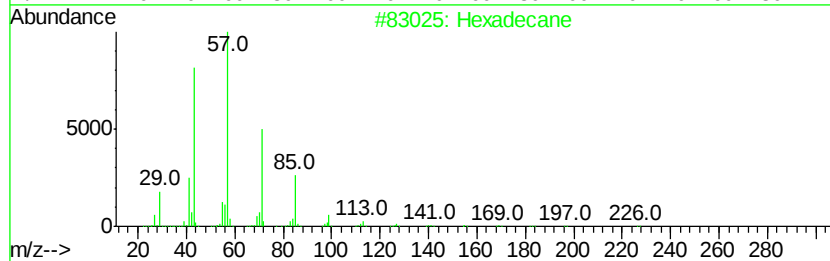
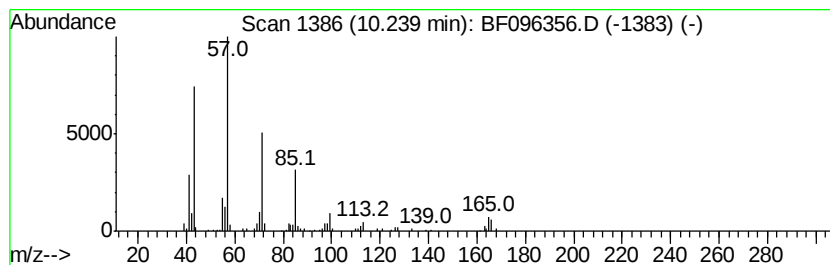
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.24	2.54 ng	122490	Acenaphthene-d10	9.79	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	86
2	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78
3	Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
4	Undecane	156	C11H24	001120-21-4	64
5	Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063017\
Data File : BF096356.D
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ClientSampleId :
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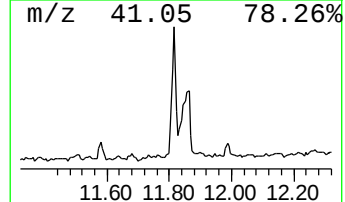
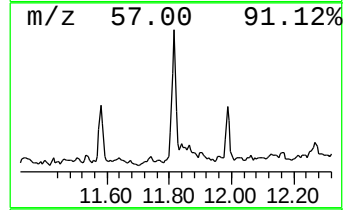
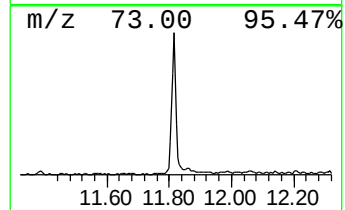
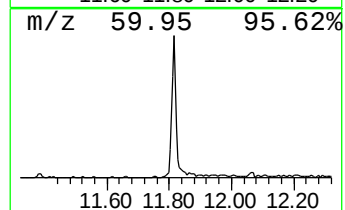
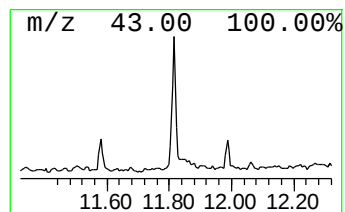
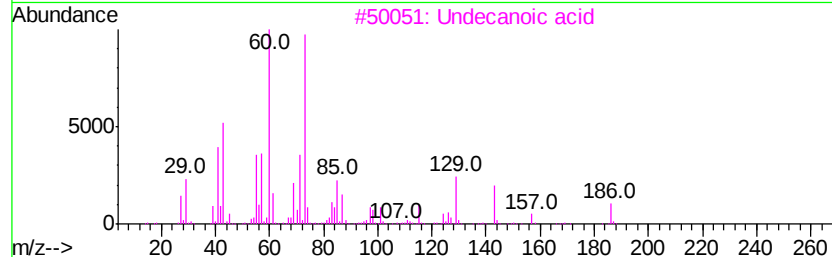
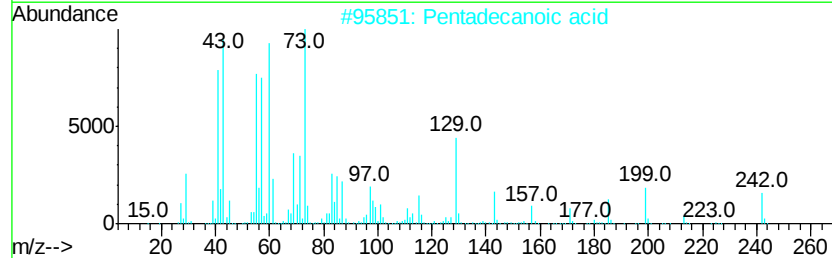
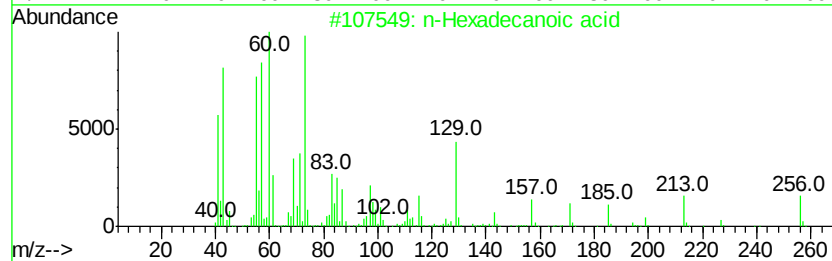
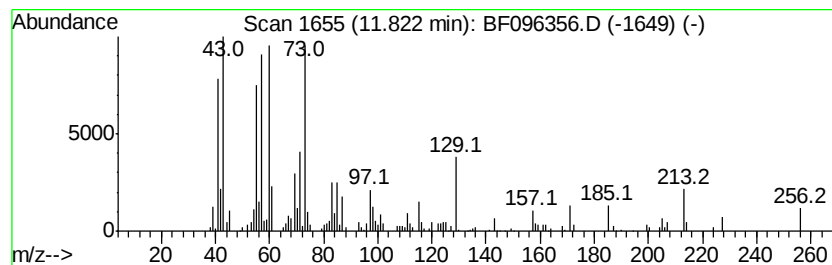
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	11.88 ng	477509	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Undecanoic acid	186	C11H22O2	000112-37-8	86
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5			Tridecanoic acid	214	C13H26O2	000638-53-9	81



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063017\
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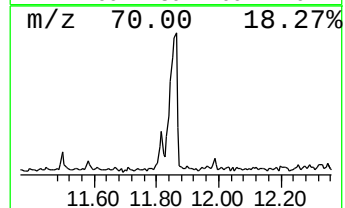
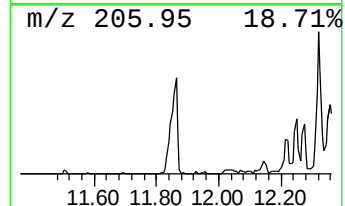
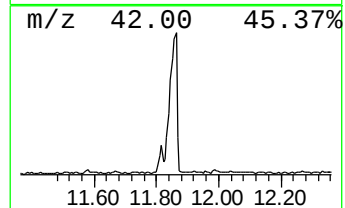
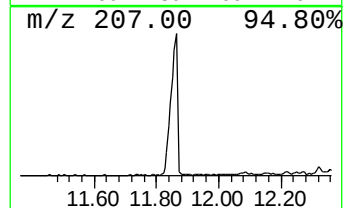
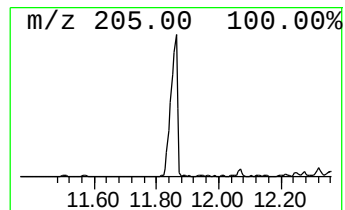
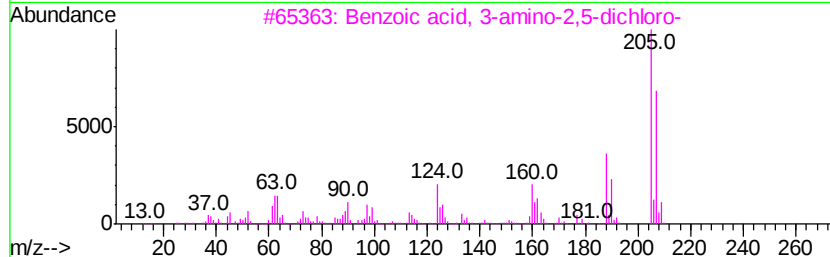
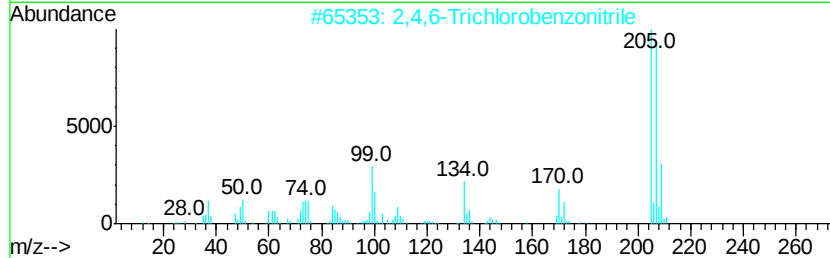
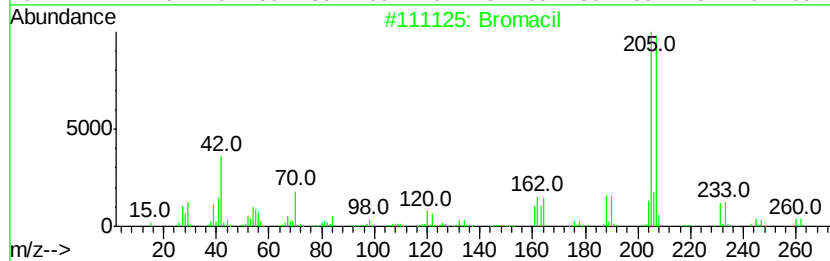
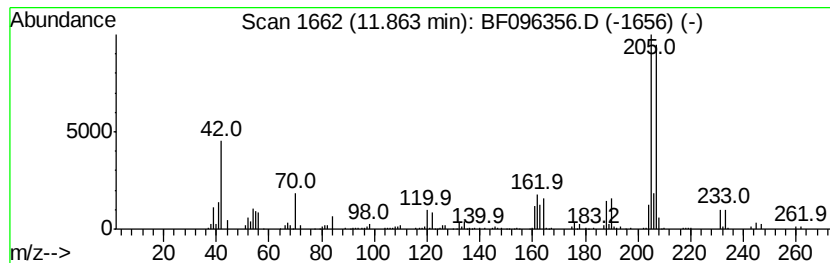
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Bromacil Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	19.66 ng	789795	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromacil	260	C9H13BrN2O2	000314-40-9	90
2		2,4,6-Trichlorobenzonitrile	205	C7H2Cl3N	006575-05-9	50
3		Benzoic acid, 3-amino-2,5-dichloro-	205	C7H5Cl2NO2	000133-90-4	40
4		5-Acetamido-2,4-dichlorobenzoic ...	247	C9H7Cl2NO3	050602-49-8	40
5		7-Chloro-2,3-dihydrofuro(2,3-b)q...	205	C11H8ClNO	064124-91-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
Data File : BF096356.D
Acq On : 1 Jul 2017 00:32
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ClientSampleId :
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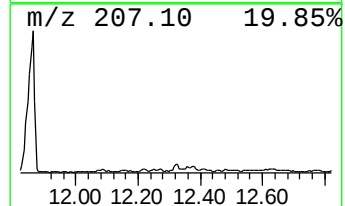
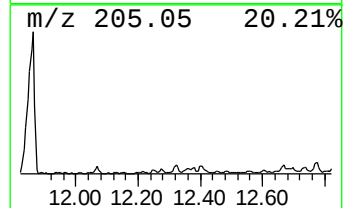
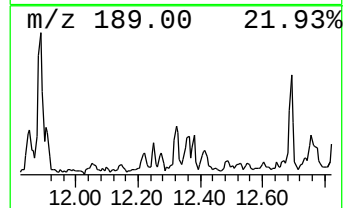
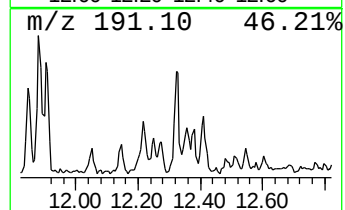
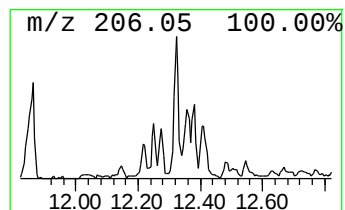
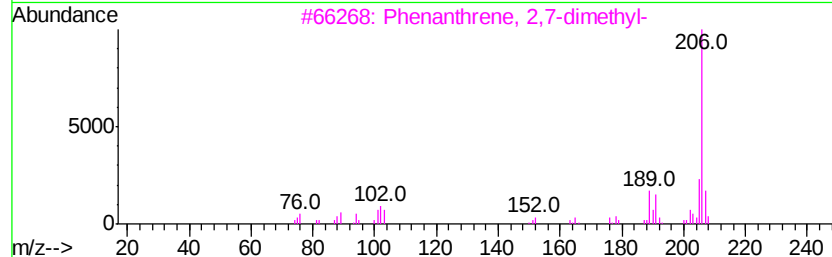
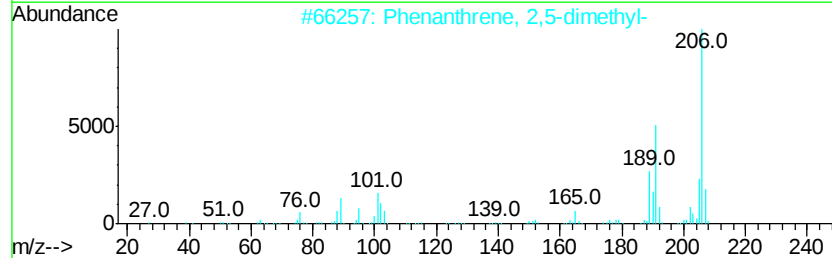
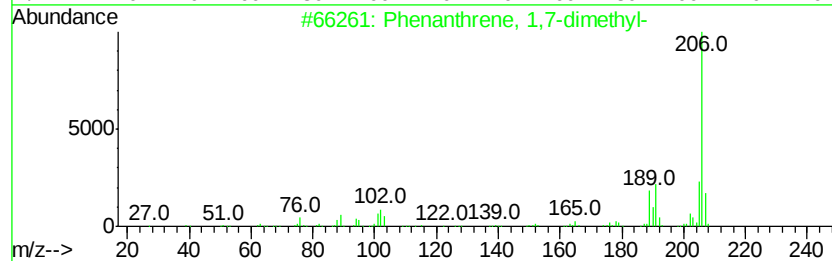
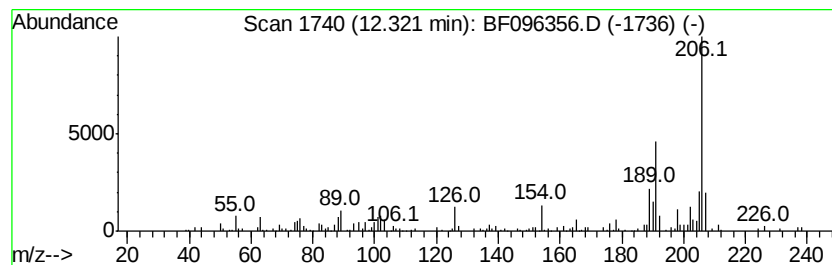
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenanthrene, 1,7-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	3.03 ng	121779	Phenanthrene-d10	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	91
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	90
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
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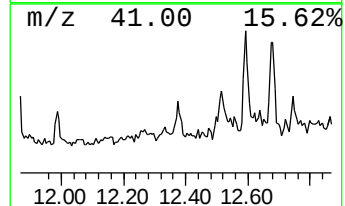
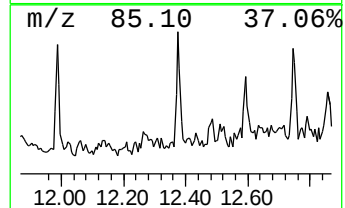
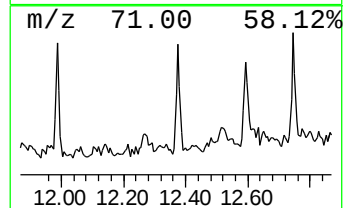
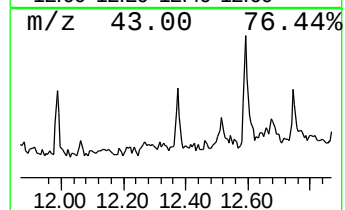
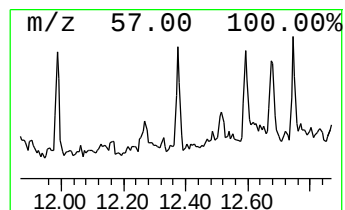
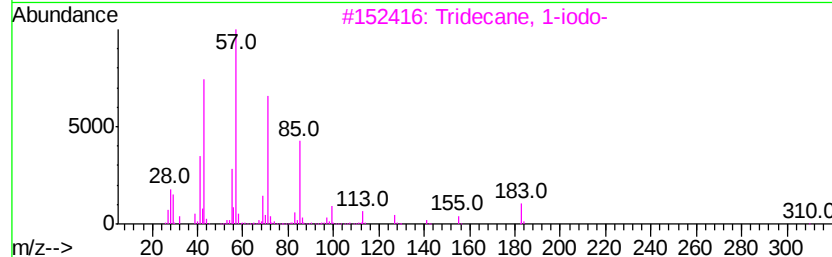
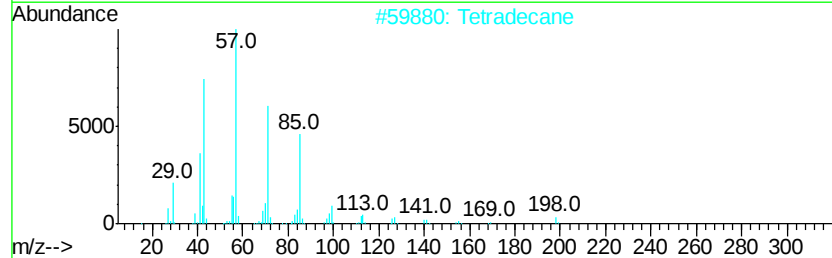
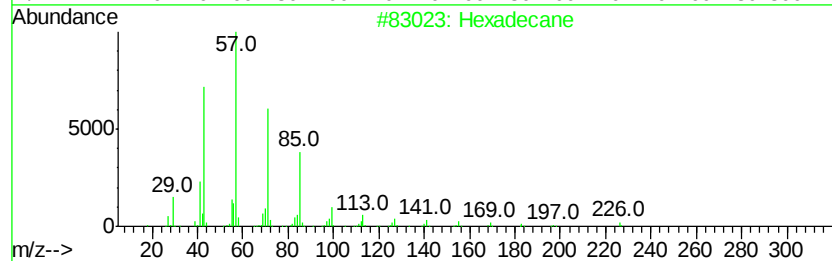
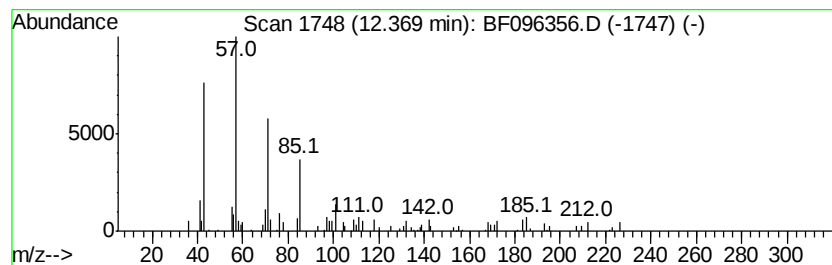
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Tetratetracontane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	2.18 ng	87479	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	87
2		Tetradecane	198	C14H30	000629-59-4	80
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
4		Tetratetracontane	619	C44H90	007098-22-8	72
5		Docosane	310	C22H46	000629-97-0	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
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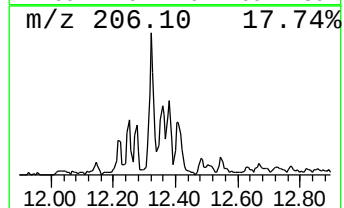
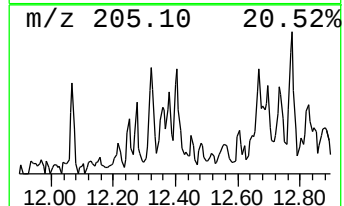
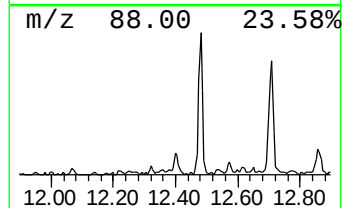
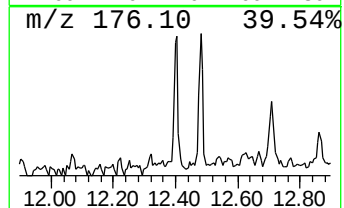
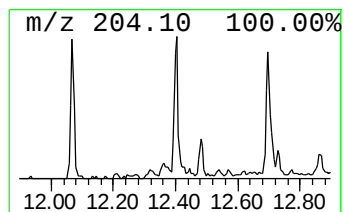
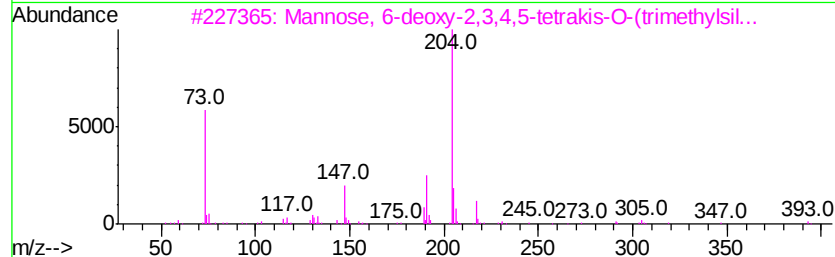
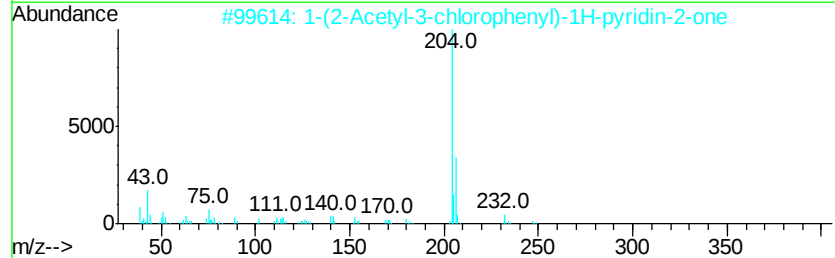
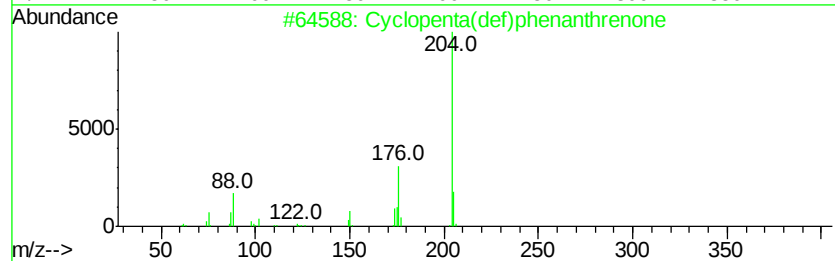
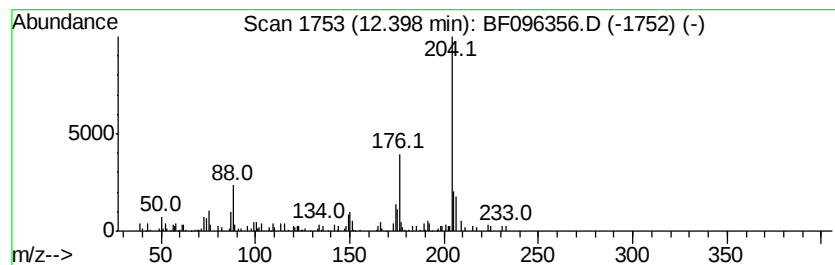
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclopenta(def)phenanthrenone Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	2.08 ng	83723	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	95
2		1-(2-Acetyl-3-chlorophenyl)-1H-p...	247	C13H10ClNO2	1000306-70-9	47
3		Mannose, 6-deoxy-2,3,4,5-tetraki...	452	C18H44O5Si4	019127-15-2	47
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	46
5		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063017\
Data File : BF096356.D
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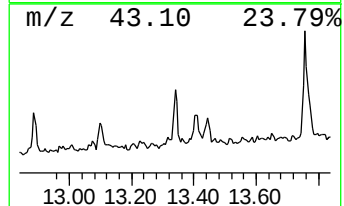
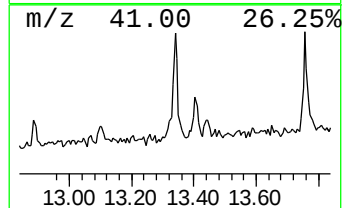
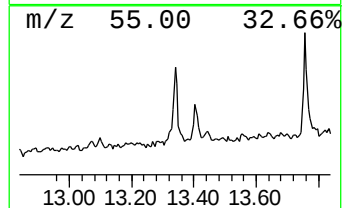
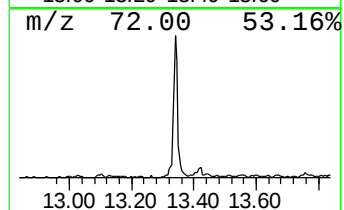
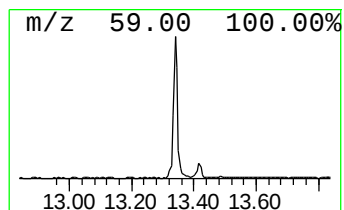
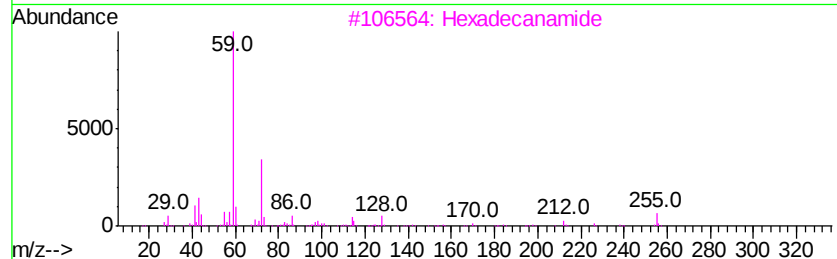
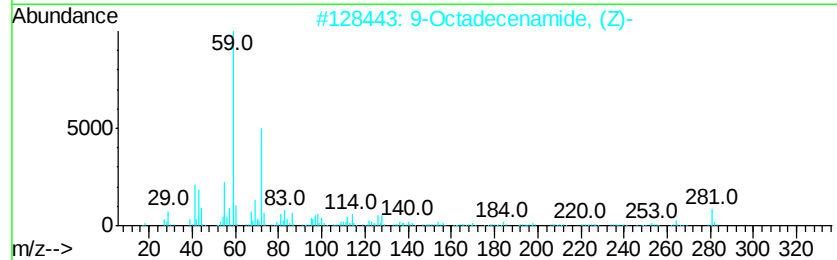
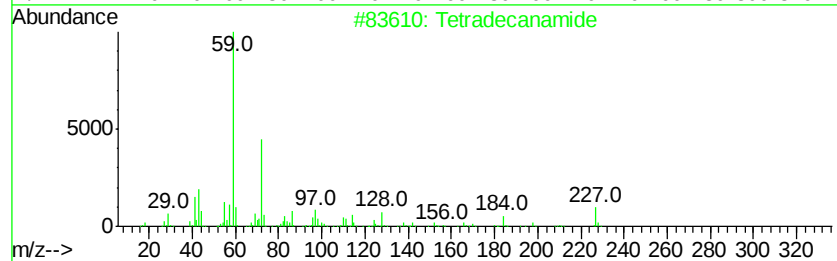
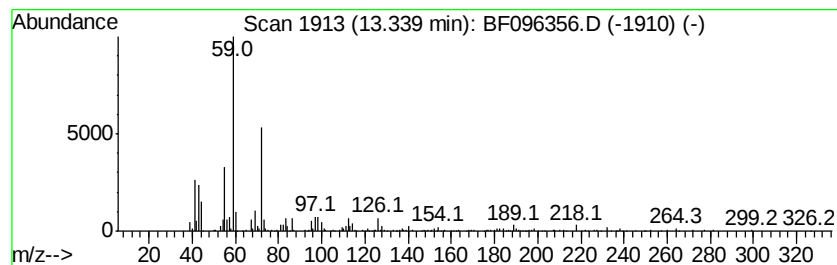
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetradecanamide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	4.54 ng	334895	Chrysene-d12	13.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanamide	227	C14H29NO	000638-58-4	87
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3			Hexadecanamide	255	C16H33NO	000629-54-9	72
4			Nonanamide	157	C9H19NO	001120-07-6	59
5			Dodecanamide	199	C12H25NO	001120-16-7	45



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063017\
Data File : BF096356.D
Acq On : 1 Jul 2017 00:32
Operator : SJ/MA
Sample : I3930-03
Misc :
ALS Vial : 33 Sample Multiplier: 1

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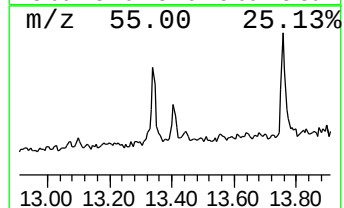
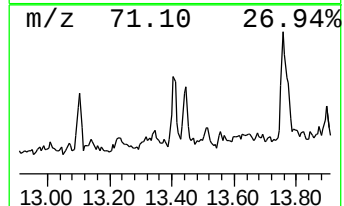
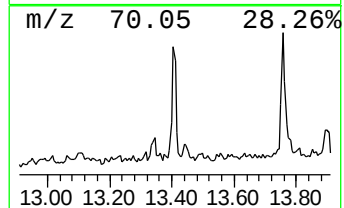
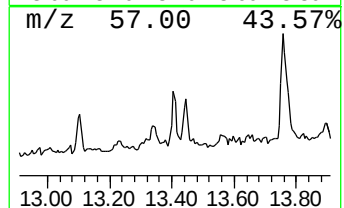
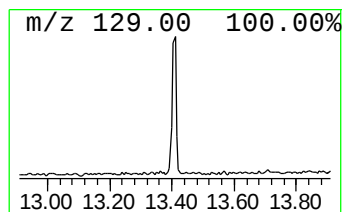
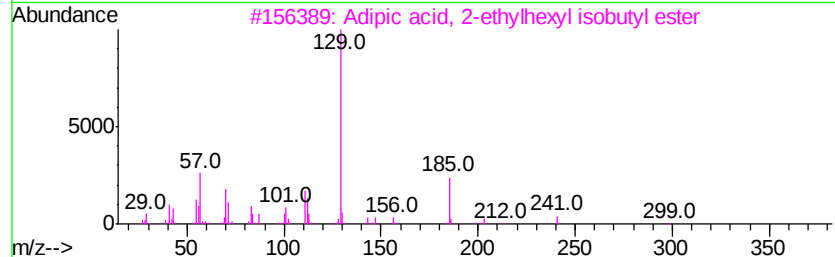
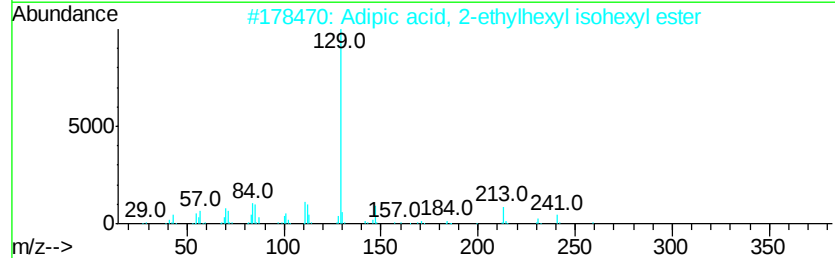
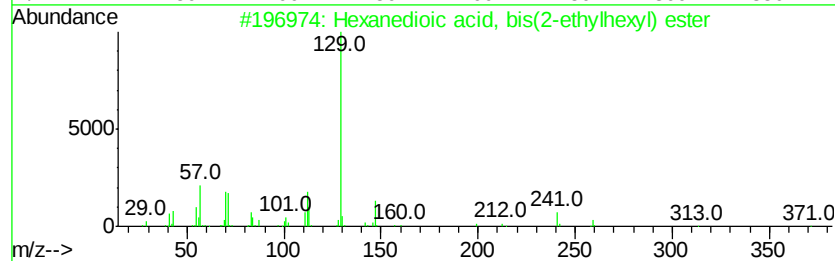
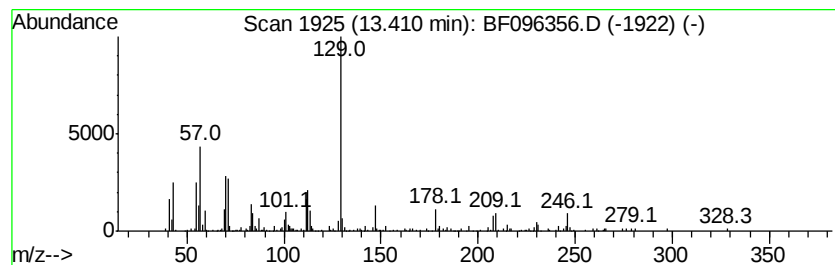
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexanedioic acid, bis(2-eth... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	2.40 ng	176722	Chrysene-d12	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	68
2		Adipic acid, 2-ethylhexyl isohex...	342	C20H38O4	1000324-48-3	58
3		Adipic acid, 2-ethylhexyl isobut...	314	C18H34O4	1000324-48-0	50
4		Adipic acid, isohexyl trans-2-me...	326	C19H34O4	1000324-74-8	50
5		Adipic acid, dicyclopentylmethyl...	310	C18H30O4	1000324-78-8	49



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Operator : SJ/MA
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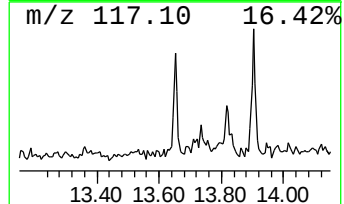
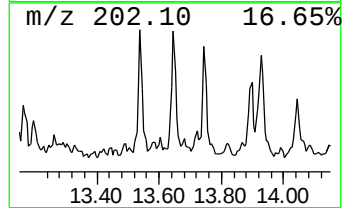
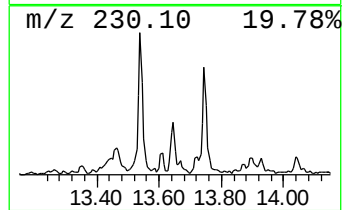
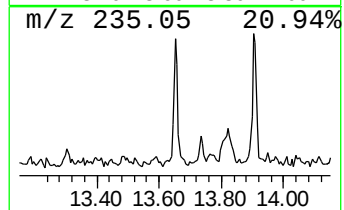
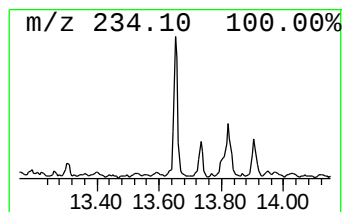
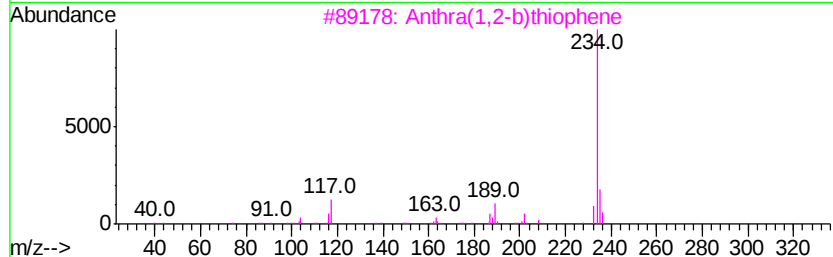
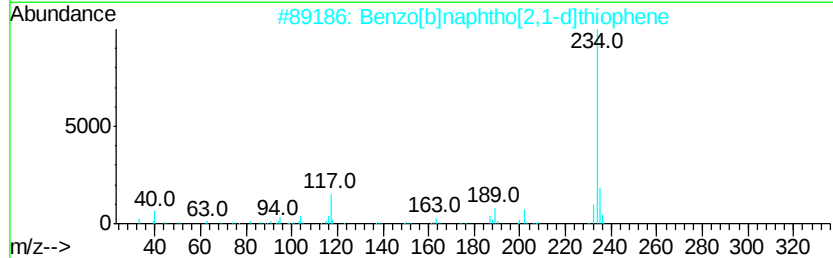
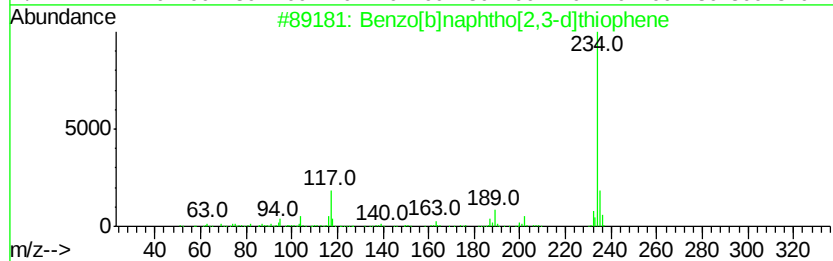
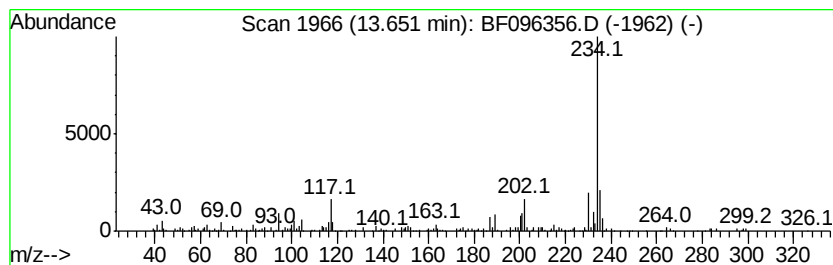
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[b]naphtho[2,3-d]thiop... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	2.38 ng	175672	Chrysene-d12	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	96
2		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	92
3		Anthra(1,2-b)thiophene	234	C16H10S	000227-86-1	86
4		Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	70
5		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
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Operator : SJ/MA
Sample : I3930-03
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B2-4-6

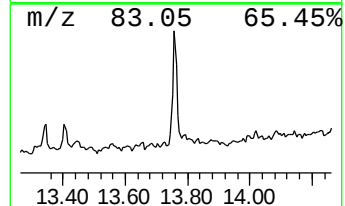
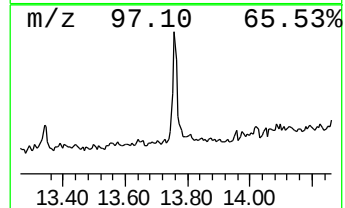
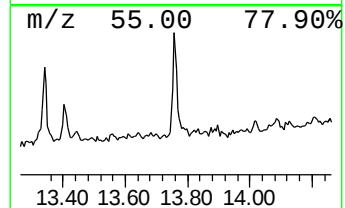
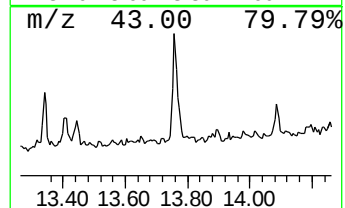
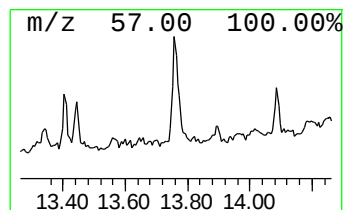
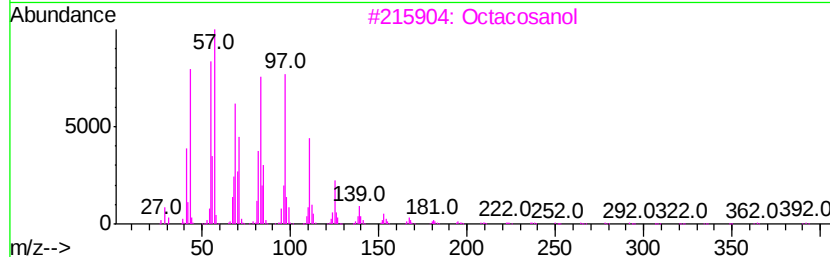
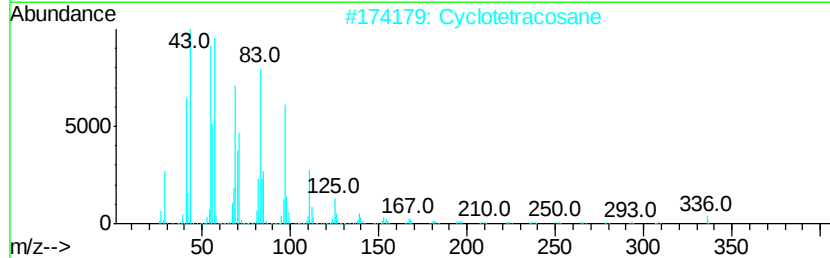
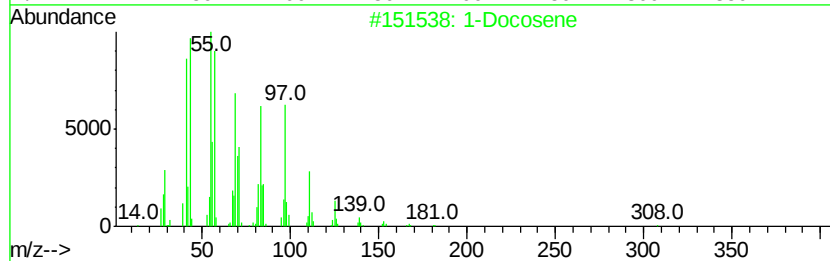
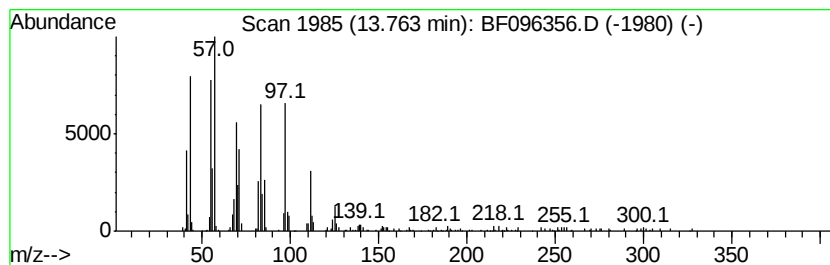
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 1-Docosene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	5.00 ng	368310	Chrysene-d12	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	94
2		Cyclotetracosane	336	C24H48	000297-03-0	94
3		Octacosanol	410	C28H58O	000557-61-9	90
4		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	90
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	90



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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B2-4-6

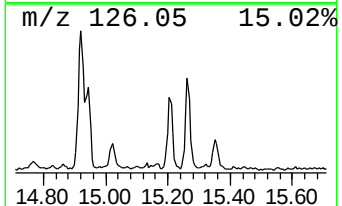
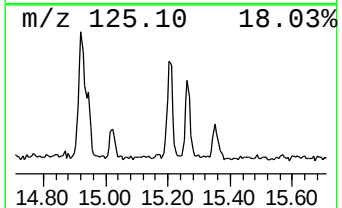
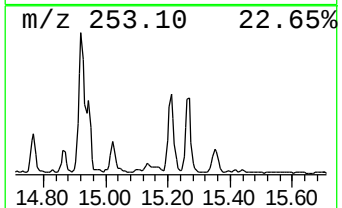
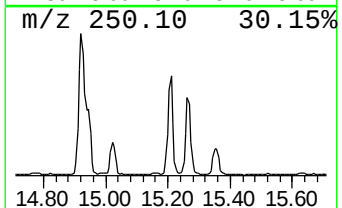
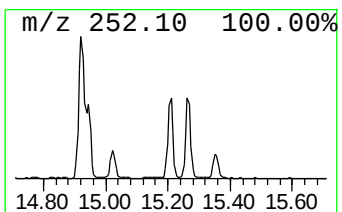
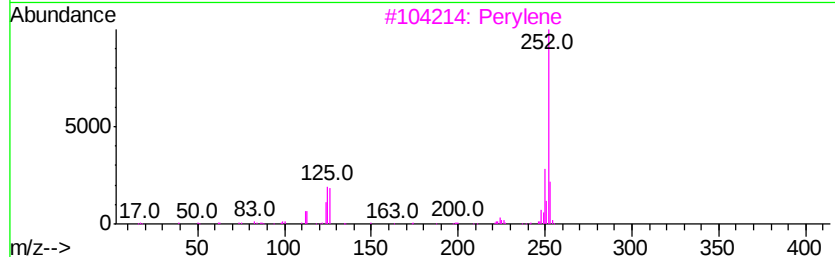
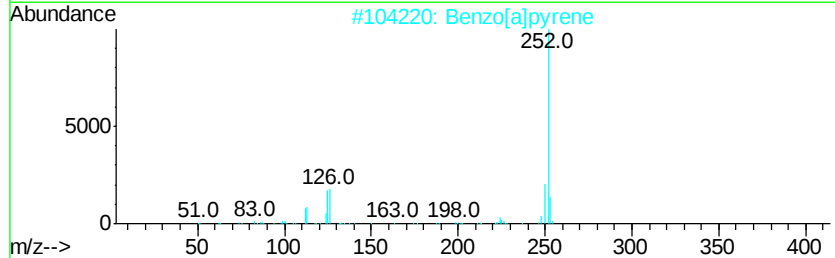
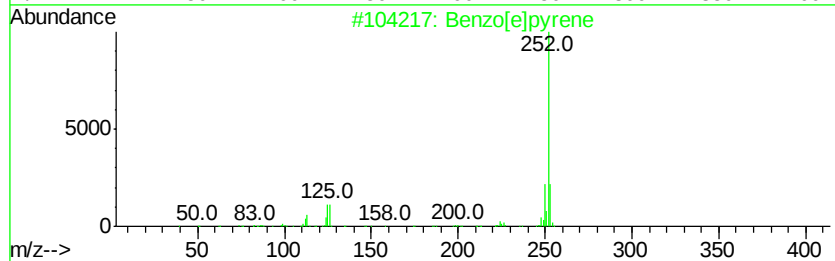
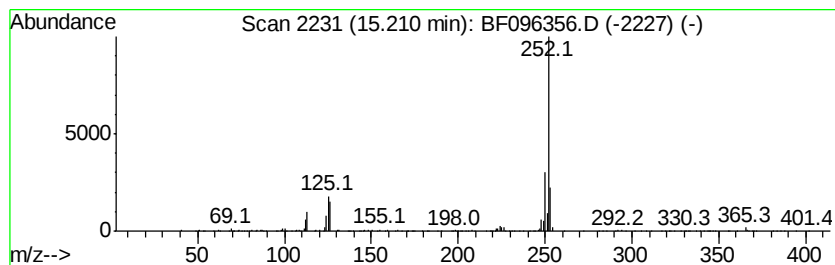
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	18.44 ng	442810	Perylene-d12	15.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Perylene	252	C20H12	000198-55-0	95
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063017\
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 Acq On : 1 Jul 2017 00:32
 Operator : SJ/MA
 Sample : I3930-03
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B2-4-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.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...	4.96	16.1 ng		694194	1	6.76	860858	20.0
unknown6.53	6.53	86.8 ng		3735420	1	6.76	860858	20.0
Tridecane	9.74	2.2 ng		107318	3	9.79	965238	20.0
Hexadecane	10.24	2.5 ng		122490	3	9.79	965238	20.0
n-Hexadecanoic acid	11.82	11.9 ng		477509	4	11.27	803660	20.0
Bromacil	11.86	19.7 ng		789795	4	11.27	803660	20.0
Phenanthrene, 1,7...	12.32	3.0 ng		121779	4	11.27	803660	20.0
Tetratetracontane	12.37	2.2 ng		87479	4	11.27	803660	20.0
Cyclopenta(def)ph...	12.40	2.1 ng		83723	4	11.27	803660	20.0
Tetradecanamide	13.34	4.5 ng		334895	5	13.90	1474320	20.0
Hexanedioic acid,...	13.41	2.4 ng		176722	5	13.90	1474320	20.0
Benzo[b]naphtho[2...	13.65	2.4 ng		175672	5	13.90	1474320	20.0
1-Docosene	13.76	5.0 ng		368310	5	13.90	1474320	20.0
Benzo[e]pyrene	15.21	18.4 ng		442810	6	15.33	480375	20.0