

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
 Data File : BF076993.D
 Acq On : 21 Jan 2015 9:57
 Operator : IZ / TP
 Sample : G1111-06
 Misc : S
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBDC6-011615

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-BF011415.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	1.441	29	31	37	rBV	32966	60284	5.05%	0.400%
2	4.024	254	257	264	rBV	56918	133558	11.20%	0.887%
3	4.984	339	341	353	rBV	348439	760319	63.75%	5.050%
4	5.305	364	369	375	rVB	45533	79940	6.70%	0.531%
5	7.339	544	547	551	rBV	427808	734783	61.61%	4.881%
6	7.419	551	554	561	rVB	498106	876417	73.48%	5.821%
7	7.910	594	597	601	rBV	110301	167648	14.06%	1.114%
8	8.231	622	625	629	rBV	379006	597846	50.13%	3.971%
9	9.213	707	711	714	rBV	324632	506929	42.50%	3.367%
10	10.825	848	852	855	rBV	143665	240231	20.14%	1.596%
11	12.300	978	981	985	rVB	35525	61210	5.13%	0.407%
12	13.477	1080	1084	1087	rBV	590402	894437	75.00%	5.941%
13	14.540	1173	1177	1180	rBV	85429	125619	10.53%	0.834%
14	14.951	1209	1213	1216	rVB	175974	267154	22.40%	1.775%
15	15.397	1250	1252	1255	rVB	8991	14063	1.18%	0.093%
16	15.877	1290	1294	1301	rBV2	5097	13933	1.17%	0.093%
17	16.197	1319	1322	1325	rVB	10703	16195	1.36%	0.108%
18	16.586	1353	1356	1358	rBV	130664	190163	15.94%	1.263%
19	16.631	1358	1360	1363	rVV	409870	647875	54.32%	4.303%
20	17.294	1413	1418	1420	rBV4	8206	14900	1.25%	0.099%
21	17.740	1454	1457	1459	rBV	238402	286251	24.00%	1.901%
22	17.774	1459	1460	1463	rVB2	8841	11955	1.00%	0.079%
23	18.209	1496	1498	1500	rVB	53545	60931	5.11%	0.405%
24	18.369	1510	1512	1515	rBV	220969	224709	18.84%	1.493%
25	18.540	1525	1527	1530	rVB	108454	139534	11.70%	0.927%
26	18.620	1530	1534	1537	rBV3	10949	20122	1.69%	0.134%
27	18.734	1541	1544	1550	rBV	137507	208244	17.46%	1.383%
28	18.917	1558	1560	1564	rBV3	13618	26371	2.21%	0.175%
29	18.997	1564	1567	1570	rBV	698015	764833	64.13%	5.080%
30	19.089	1572	1575	1576	rBV	289004	398913	33.45%	2.650%
31	19.192	1582	1584	1586	rVB	294912	296017	24.82%	1.966%
32	19.260	1586	1590	1592	rVV	334571	530073	44.44%	3.521%
33	19.294	1592	1593	1595	rVB	47934	39163	3.28%	0.260%
34	19.420	1602	1604	1605	rBV	29597	48125	4.04%	0.320%

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35	19.454	1605	1607	1609	rVB	1113926	1192660	100.00%	7.922%
36	19.615	1618	1621	1622	rBV3	8765	14273	1.20%	0.095%
37	19.672	1623	1626	1628	rVV	692827	866463	72.65%	5.755%
38	19.706	1628	1629	1631	rVB	23590	19757	1.66%	0.131%
39	19.752	1631	1633	1636	rBV	69805	98128	8.23%	0.652%
40	19.809	1636	1638	1640	rVB	26737	31370	2.63%	0.208%
41	19.969	1649	1652	1655	rVB	796177	834564	69.98%	5.543%
42	20.517	1696	1700	1707	rBV9	7924	28350	2.38%	0.188%
43	20.655	1709	1712	1715	rVB4	52174	74058	6.21%	0.492%
44	20.872	1729	1731	1733	rVB3	11629	15336	1.29%	0.102%
45	20.975	1739	1740	1743	rVB	28587	39919	3.35%	0.265%
46	21.238	1758	1763	1769	rBV	290357	478011	40.08%	3.175%
47	21.375	1772	1775	1776	rBV	24553	40800	3.42%	0.271%
48	21.443	1778	1781	1784	rBV	66407	89988	7.55%	0.598%
49	21.500	1784	1786	1788	rBV3	27314	38833	3.26%	0.258%
50	21.546	1788	1790	1793	rBV4	10303	22269	1.87%	0.148%
51	21.626	1794	1797	1799	rBV3	13261	26736	2.24%	0.178%
52	21.718	1802	1805	1808	rVV4	33509	74858	6.28%	0.497%
53	21.763	1808	1809	1813	rVV	68178	124510	10.44%	0.827%
54	21.866	1816	1818	1821	rVB4	31872	40278	3.38%	0.268%
55	21.969	1825	1827	1828	rBV2	16431	20901	1.75%	0.139%
56	22.003	1828	1830	1832	rBV	50949	73347	6.15%	0.487%
57	22.072	1832	1836	1838	rVV2	45782	93265	7.82%	0.620%
58	22.220	1847	1849	1851	rBV2	21343	31446	2.64%	0.209%
59	22.255	1851	1852	1857	rVB2	50304	86811	7.28%	0.577%
60	22.518	1873	1875	1879	rVB3	113118	148462	12.45%	0.986%
61	22.632	1883	1885	1886	rBV2	33516	51327	4.30%	0.341%
62	22.666	1886	1888	1892	rVB4	26381	33185	2.78%	0.220%
63	22.746	1893	1895	1899	rBV3	33804	87068	7.30%	0.578%
64	22.849	1899	1904	1905	rBV2	118786	272298	22.83%	1.809%
65	22.883	1905	1907	1913	rVB	228933	310563	26.04%	2.063%
66	23.101	1924	1926	1928	rBV3	32911	59991	5.03%	0.398%
67	23.478	1955	1959	1962	rBV4	65632	176305	14.78%	1.171%

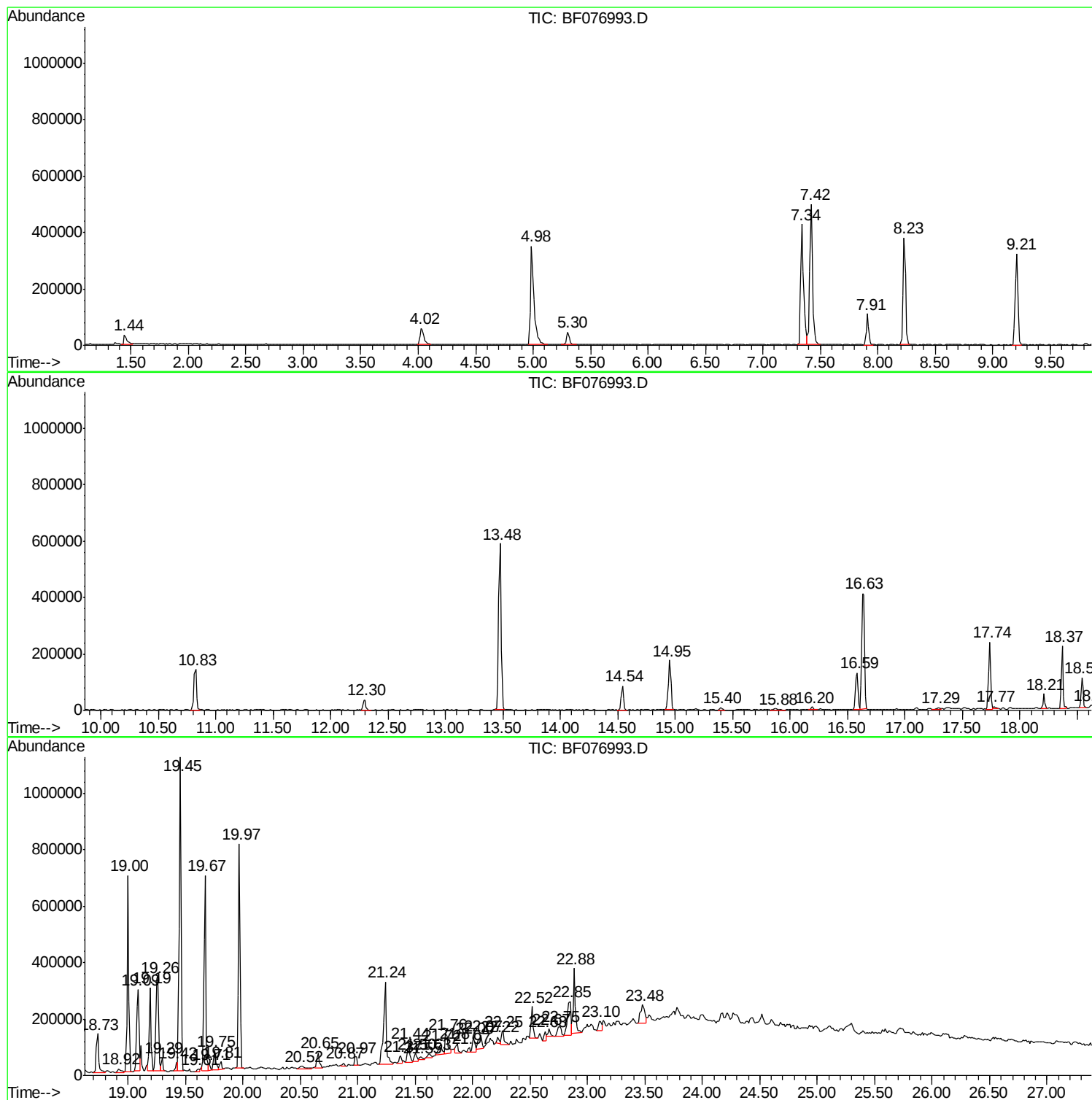
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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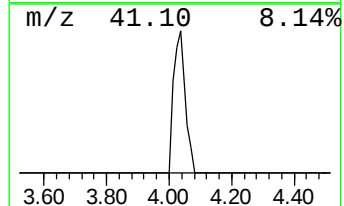
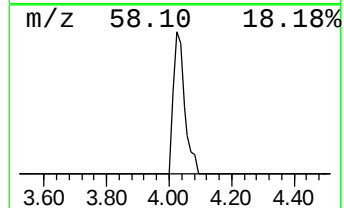
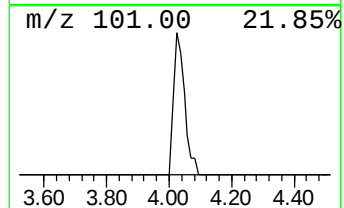
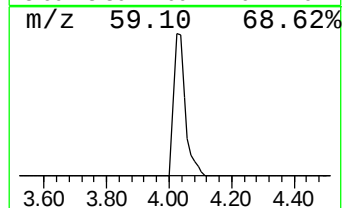
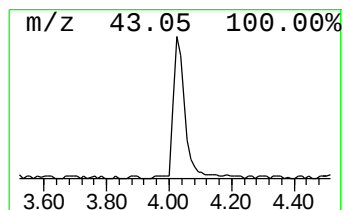
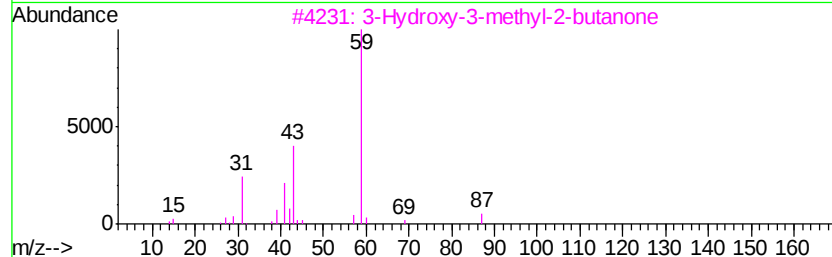
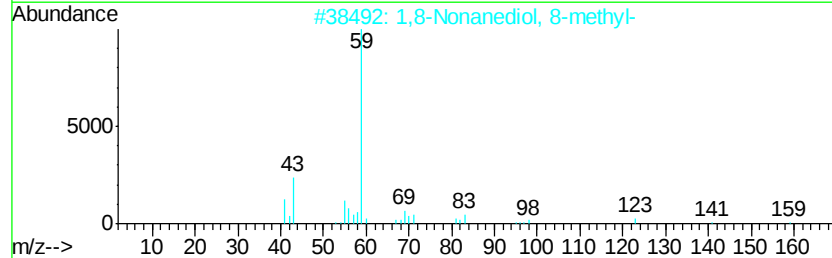
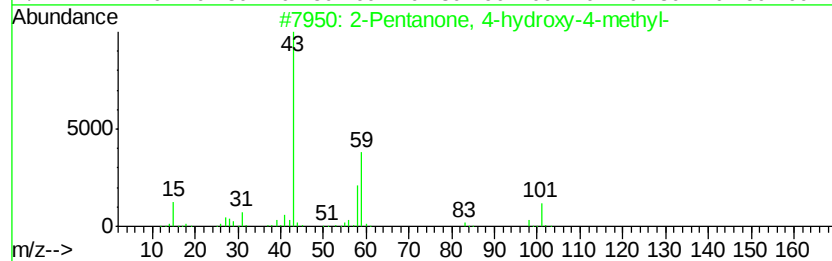
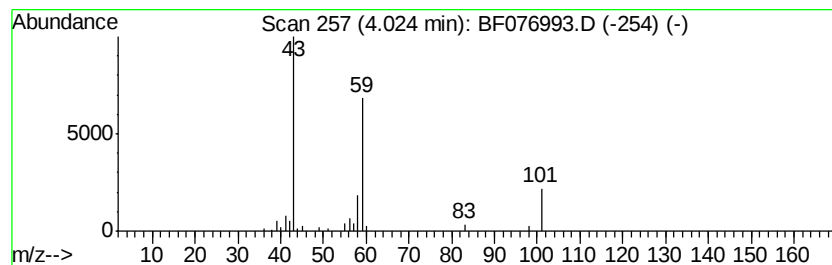
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.02	15.93 ng	133558	1,4-Dichlorobenzene-d4	7.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
2		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	32
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
4		1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	25
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23



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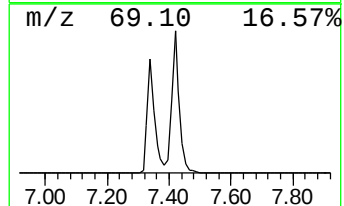
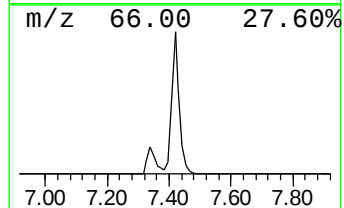
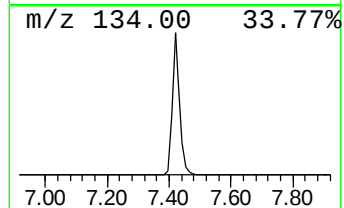
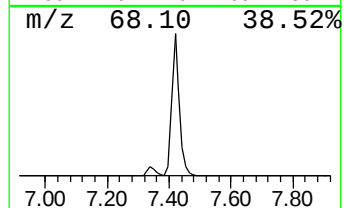
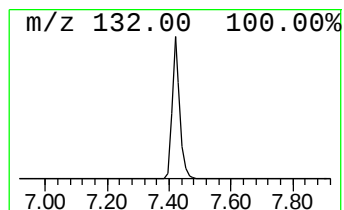
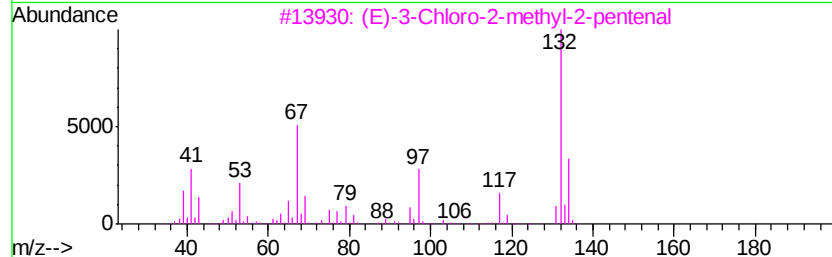
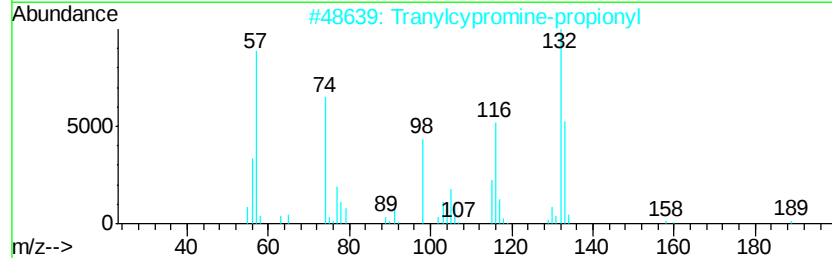
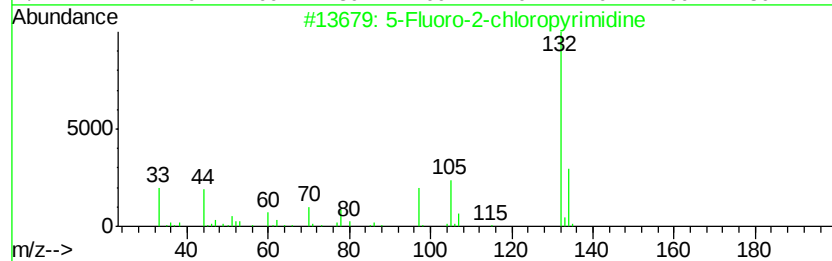
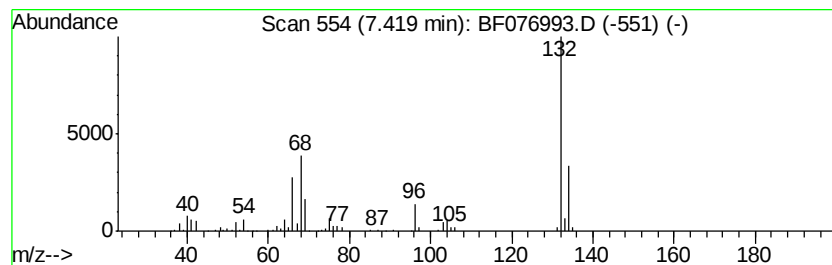
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 unknown7.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	104.55 ng	876417	1,4-Dichlorobenzene-d4	7.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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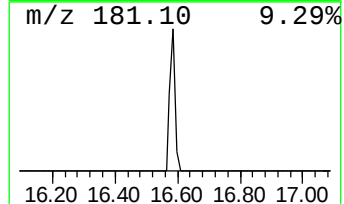
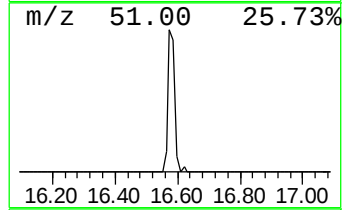
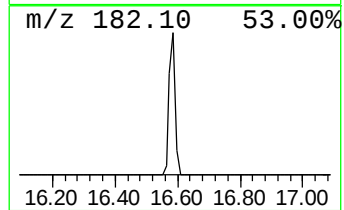
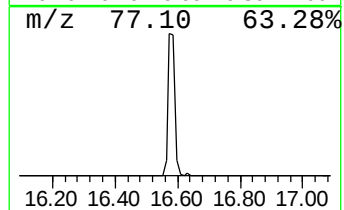
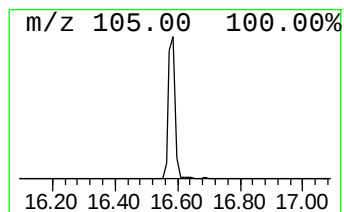
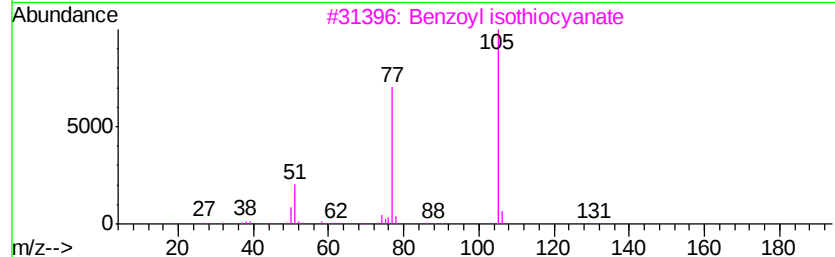
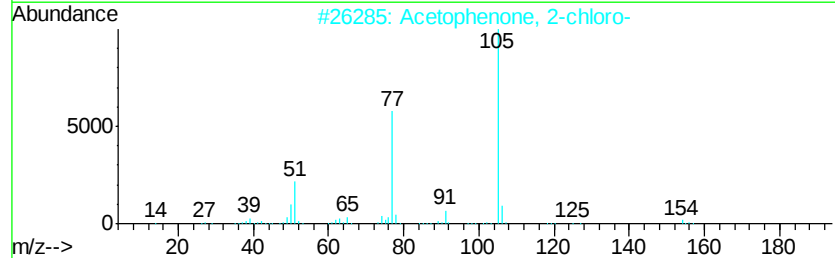
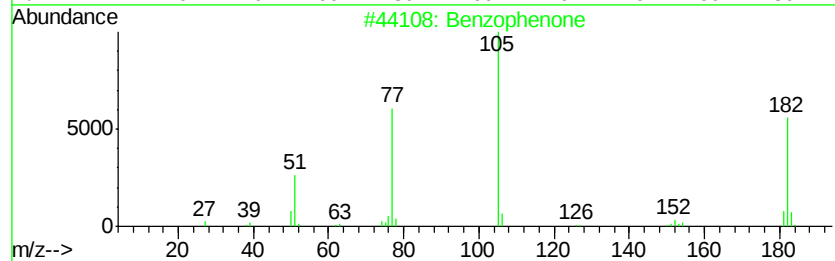
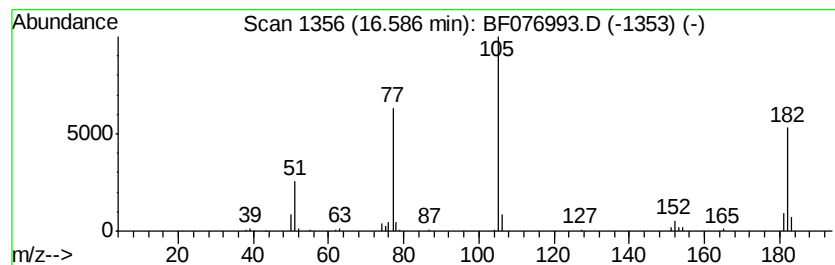
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	13.29 ng	190163	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3		Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	47
4		1-Propanone, 3-chloro-1-phenyl-	168	C9H9ClO	000936-59-4	47
5		Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	47



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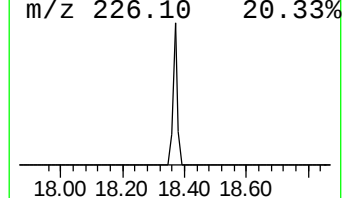
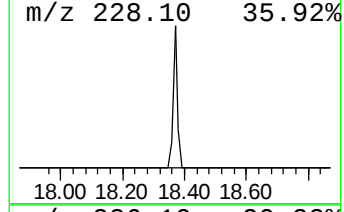
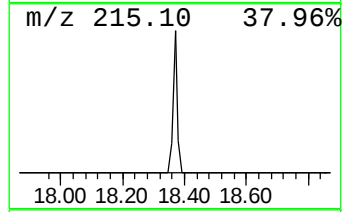
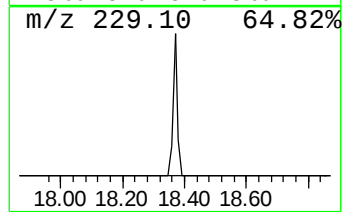
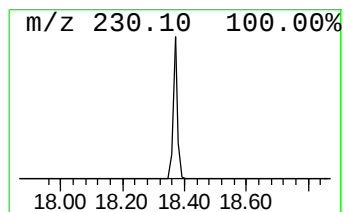
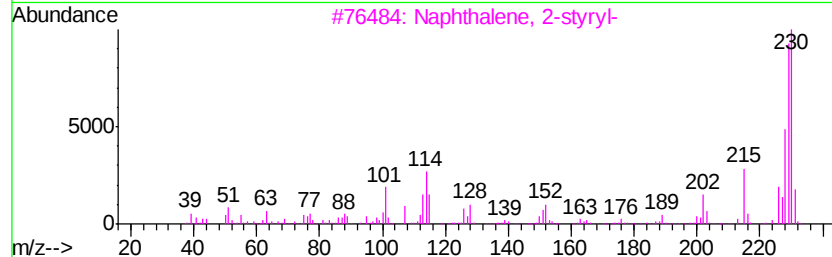
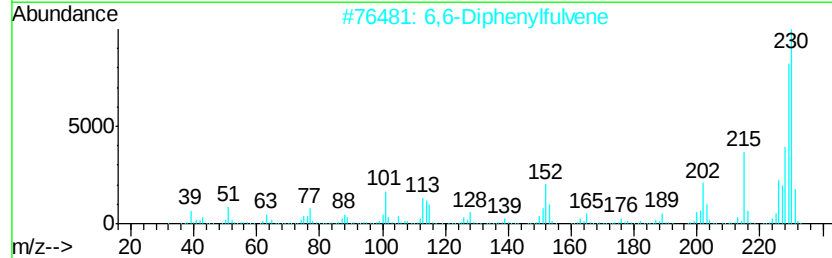
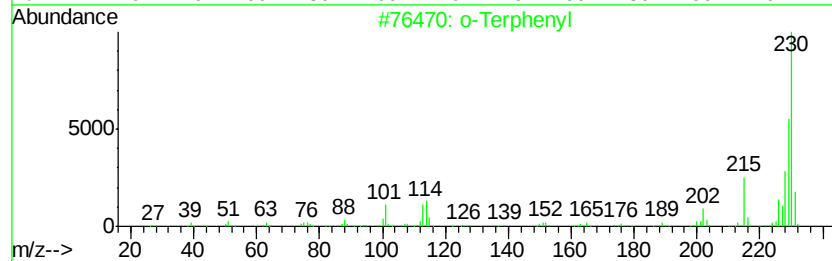
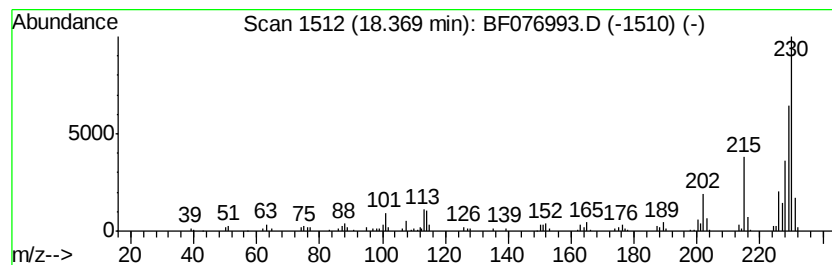
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Peak Number 7 o-Terphenyl Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	15.70 ng	224709	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Terphenyl	230	C18H14	000084-15-1	96
2		6,6-Diphenylfulvene	230	C18H14	002175-90-8	93
3		Naphthalene, 2-styryl-	230	C18H14	002039-70-5	89
4		Naphthacene, 5,12-dihydro-	230	C18H14	000959-02-4	64
5		Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	58



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Instrument :
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ClientSampleId :
RBDC6-011615

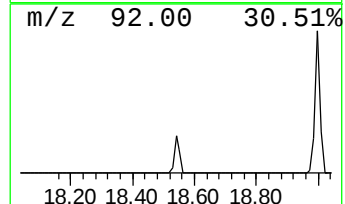
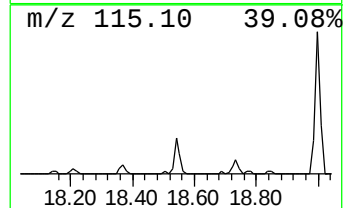
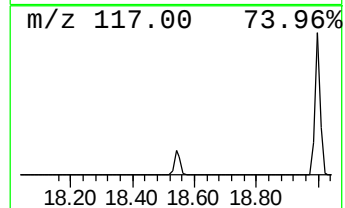
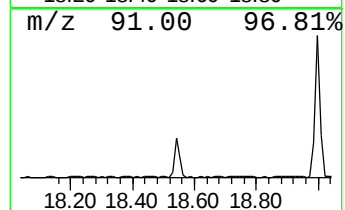
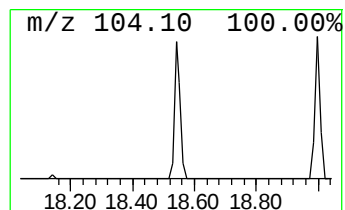
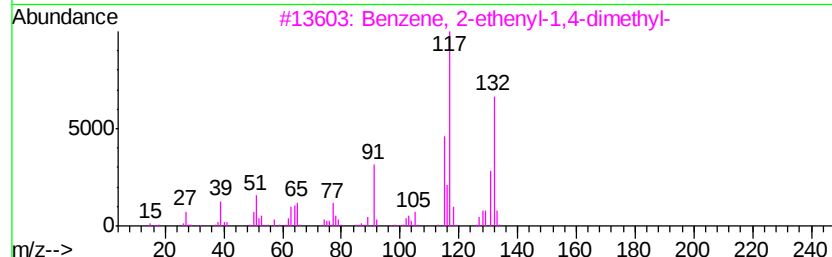
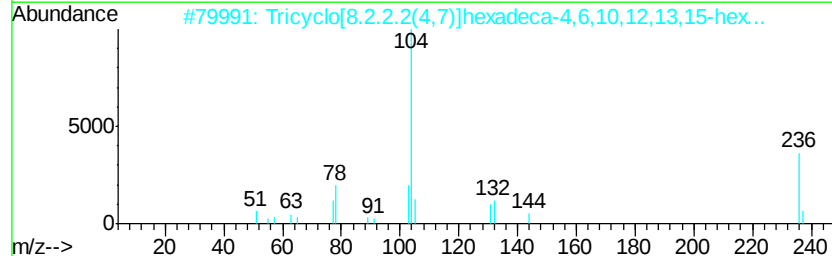
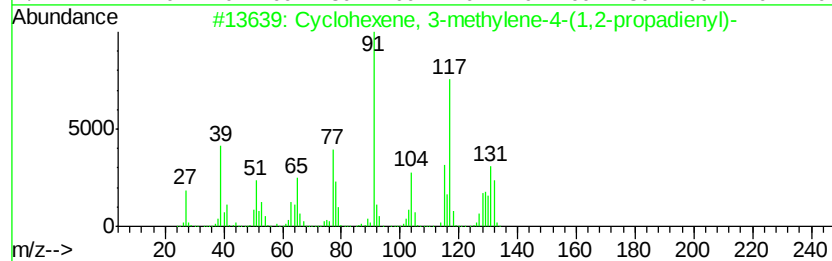
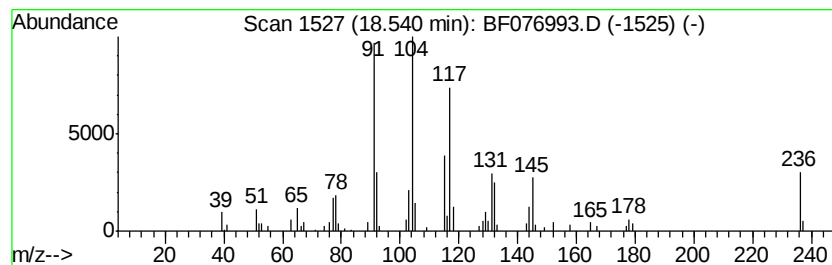
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown18.54 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	9.75 ng	139534	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 3-methylene-4-(1,2-...	132	C10H12	068377-81-1	46
2		Tricyclo[8.2.2.2(4,7)]hexadeca-4...	236	C17H16O	000729-30-6	42
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	41
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	38
5		Benzene, 1-butenyl-, (E)-	132	C10H12	001005-64-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076993.D
Acq On : 21 Jan 2015 9:57
Operator : IZ / TP
Sample : G1111-06
Misc : S
ALS Vial : 31 Sample Multiplier: 1

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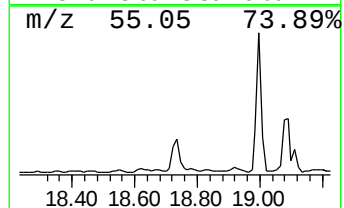
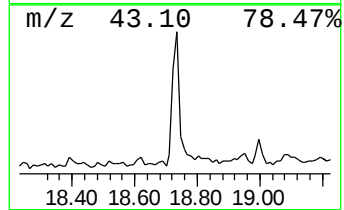
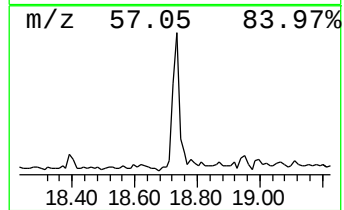
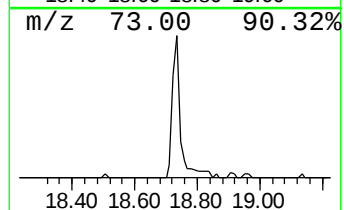
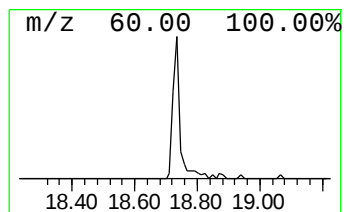
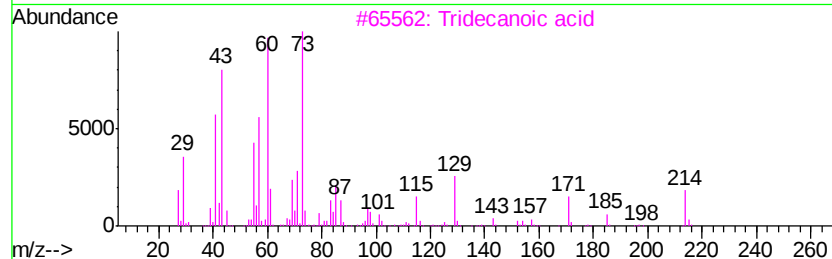
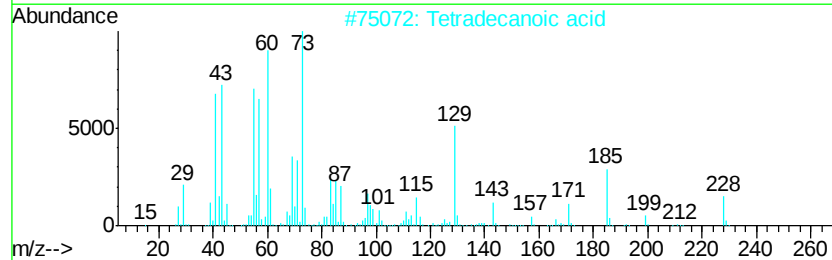
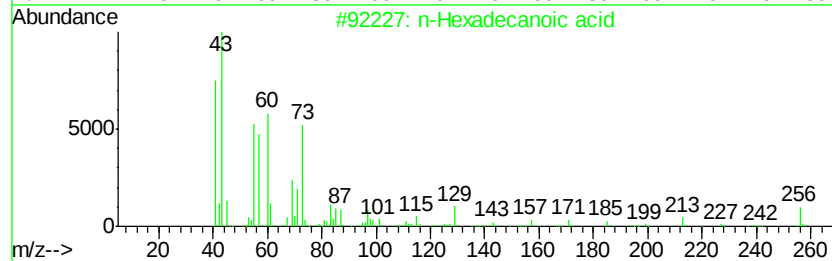
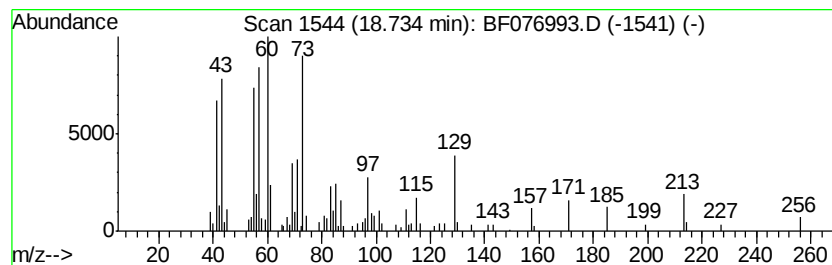
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	14.55 ng	208244	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3		Tridecanoic acid	214	C13H26O2	000638-53-9	64
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	59
5		Undecanoic acid	186	C11H22O2	000112-37-8	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076993.D
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Misc : S
ALS Vial : 31 Sample Multiplier: 1

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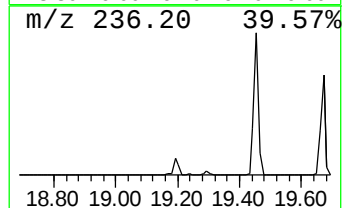
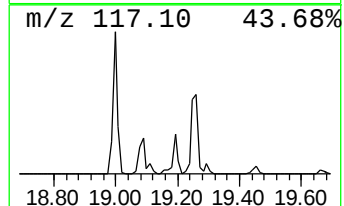
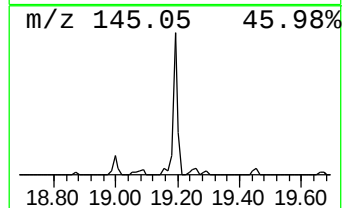
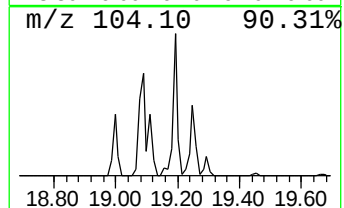
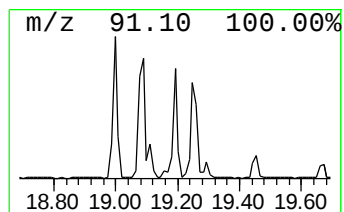
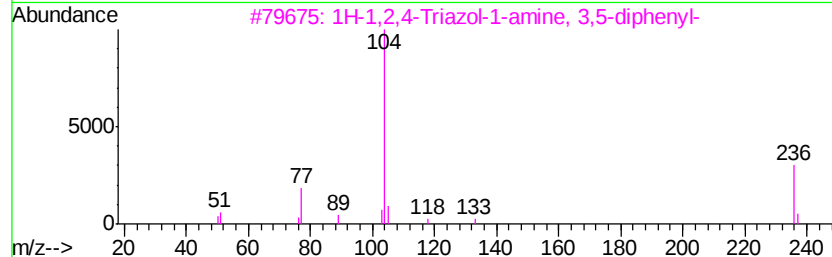
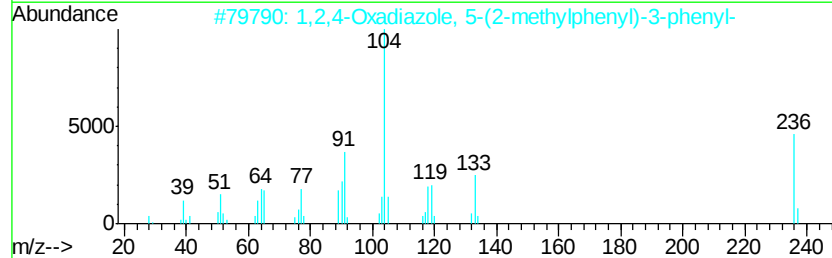
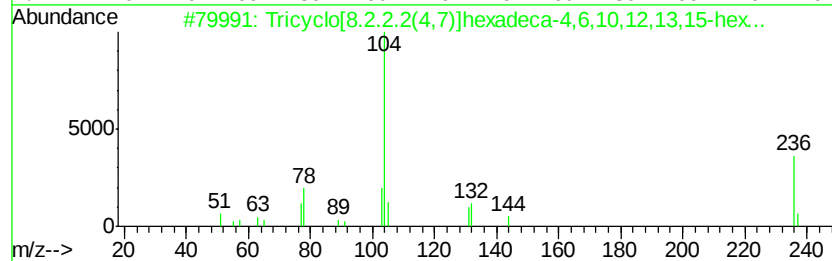
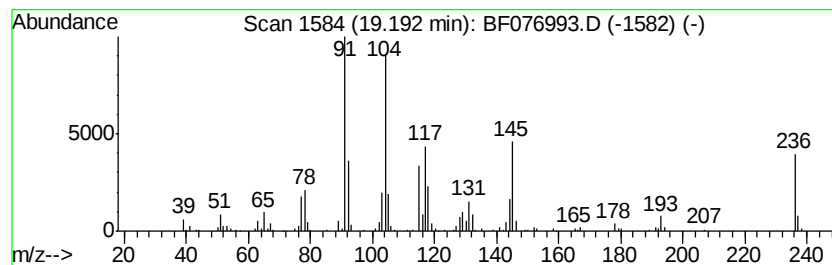
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	20.68 ng	296017	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[8.2.2.2(4,7)]hexadeca-4...	236	C17H16O	000729-30-6	87
2		1,2,4-Oxadiazole, 5-(2-methylphe...	236	C15H12N2O	054494-15-4	46
3		1H-1,2,4-Triazol-1-amine, 3,5-di...	236	C14H12N4	034985-93-8	38
4		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	30
5		3-Phenylthiolane, S,S-dioxide	196	C10H12O2S	016766-64-6	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
Data File : BF076993.D
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Misc : S
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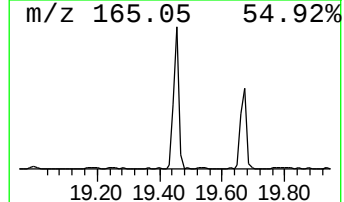
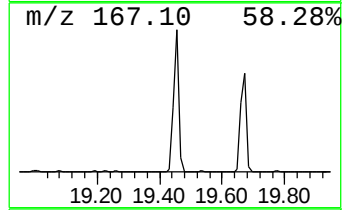
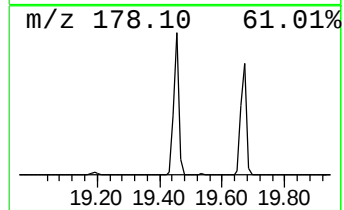
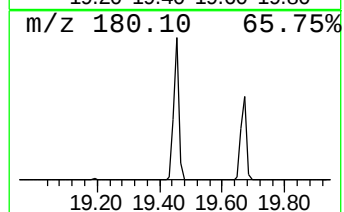
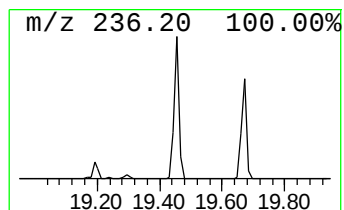
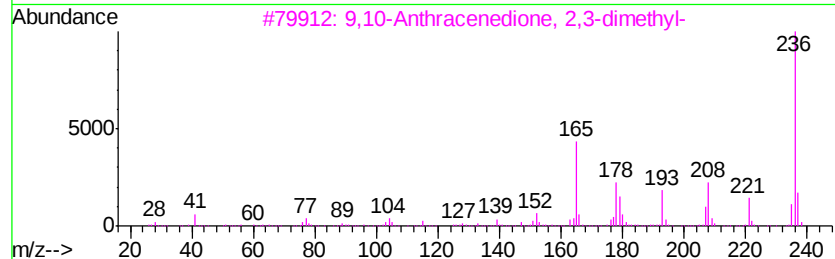
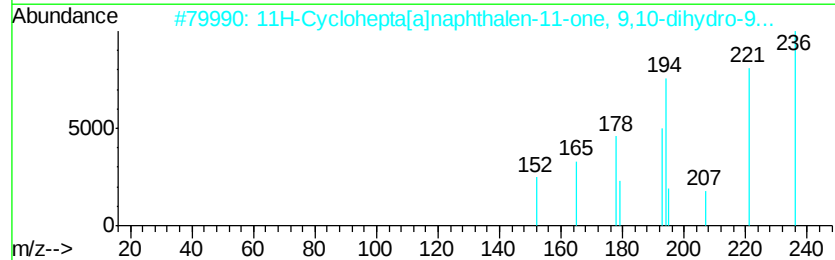
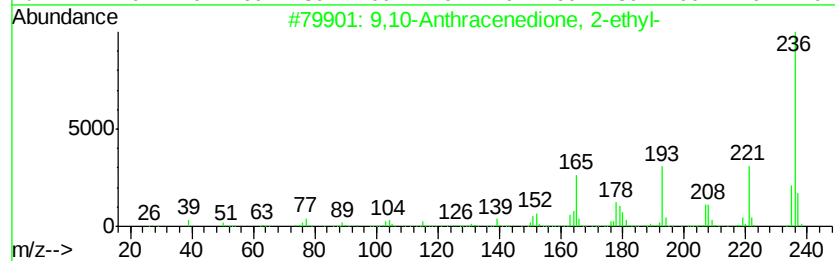
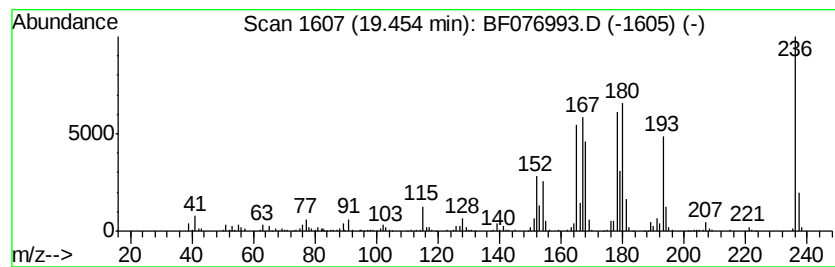
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown19.45 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.45	83.33 ng	1192660	Phenanthrene-d10	17.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione, 2-ethyl-	236	C16H12O2	000084-51-5	49
2		11H-Cyclohepta[a]naphthalen-11-one...	236	C17H16O	058111-76-5	43
3		9,10-Anthracenedione, 2,3-dimethyl-	236	C16H12O2	006531-35-7	38
4		Germacr-4-en-12-oic acid, 6.alpha...	236	C15H24O2	025861-62-5	30
5		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
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Misc : S
ALS Vial : 31 Sample Multiplier: 1

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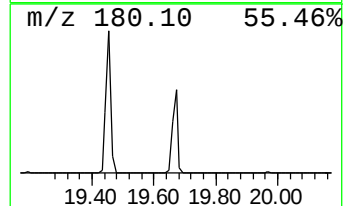
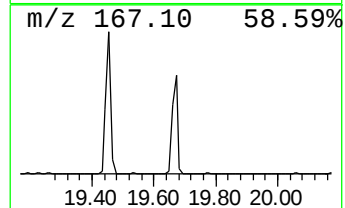
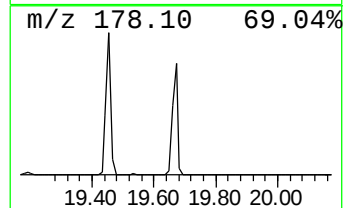
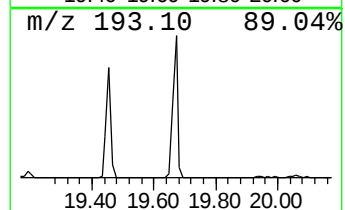
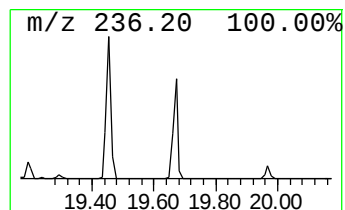
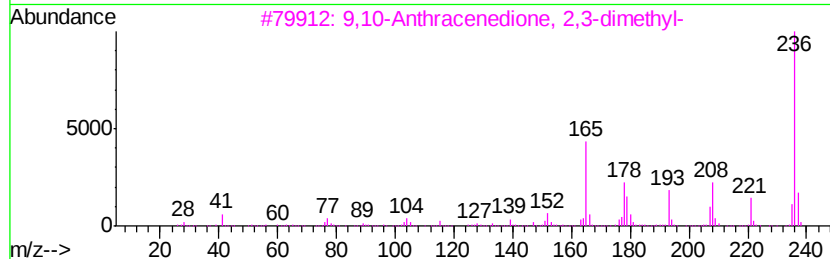
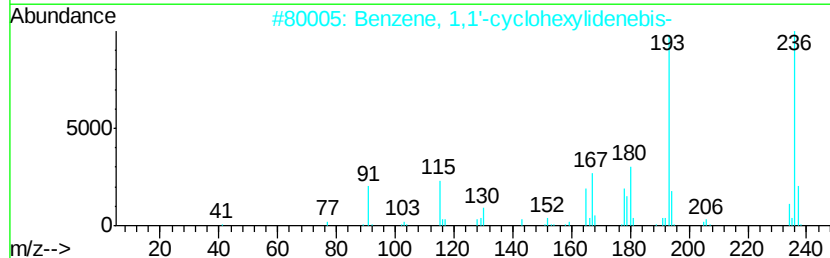
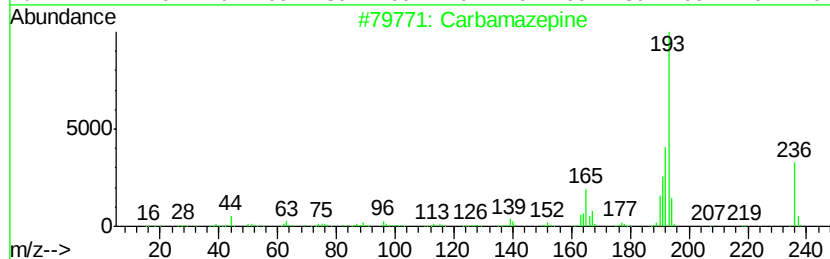
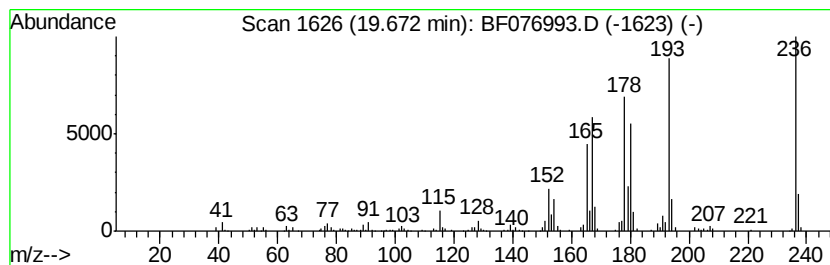
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Carbamazepine Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.67	36.25 ng	866463	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbamazepine	236	C15H12N2O	000298-46-4	60
2		Benzene, 1,1'-cyclohexylidenebis-	236	C18H20	021113-55-3	52
3		9,10-Anthracenedione, 2,3-dimethyl-	236	C16H12O2	006531-35-7	35
4		1-Formylanthraquinone	236	C15H8O3	091323-92-1	27
5		Anthracene, 9,10-dihydro-9-methy...	236	C18H20	110551-72-9	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
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ALS Vial : 31 Sample Multiplier: 1

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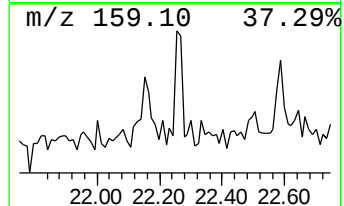
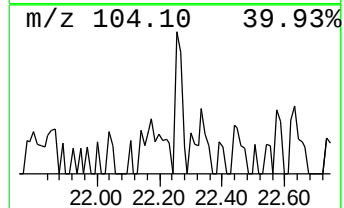
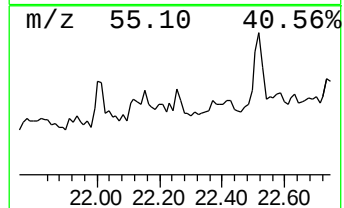
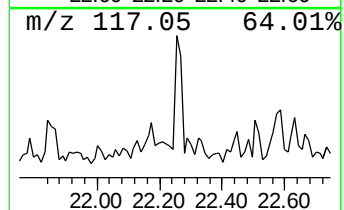
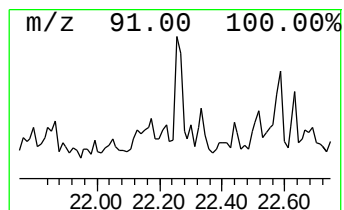
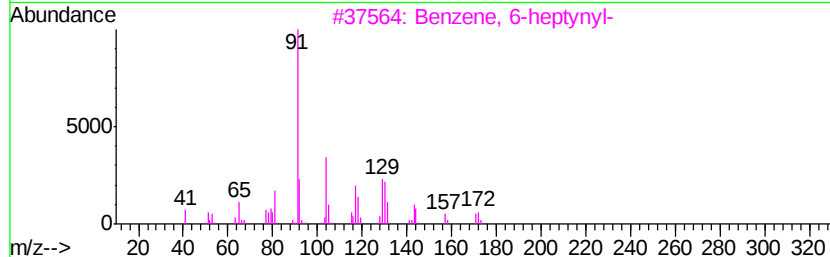
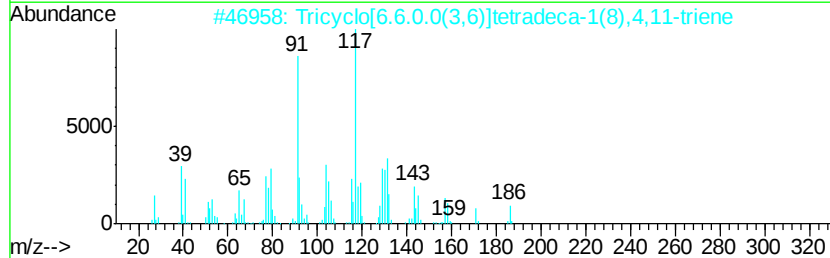
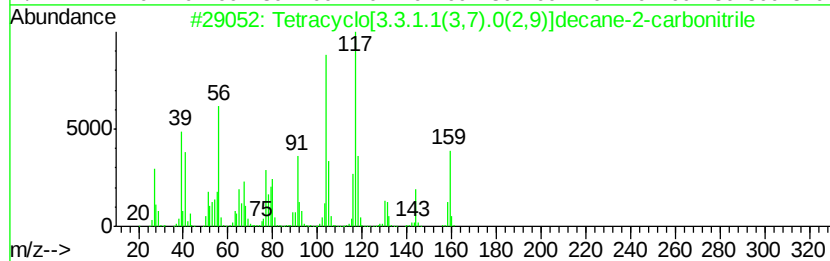
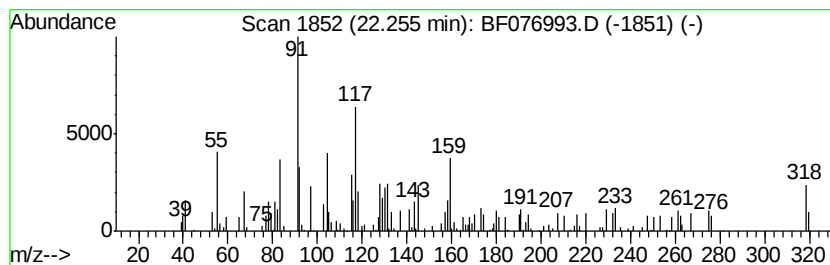
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown22.25 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.25	5.59 ng	86811	Perylene-d12	22.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracyclo[3.3.1.1(3,7).0(2,9)]d...	159	C11H13N	1000217-13-8	38
2			Tricyclo[6.6.0.0(3,6)]tetradeca...	186	C14H18	1000151-40-9	22
3			Benzene, 6-heptynyl-	172	C13H16	056293-02-8	14
4			Cyclohexene, 3-methylene-4-(1,2-...	132	C10H12	068377-81-1	10
5			2,9-Heptadecadiene-4,6-diyn-8-ol...	244	C17H24O	050816-77-8	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
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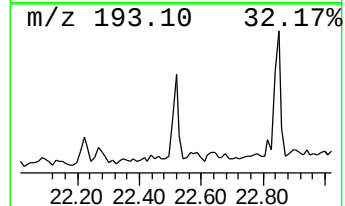
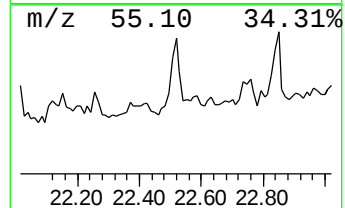
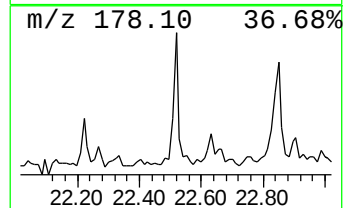
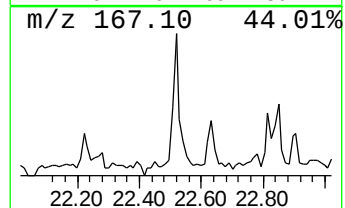
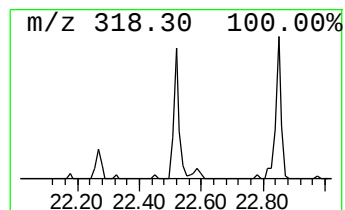
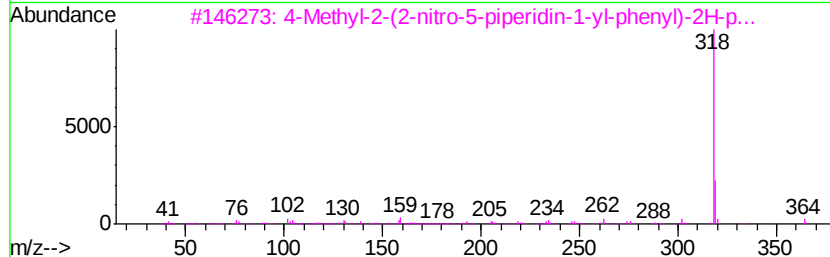
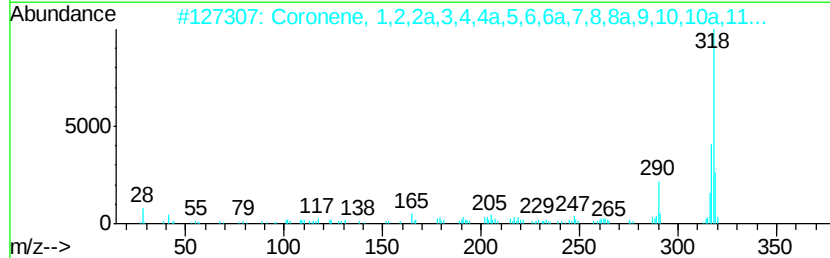
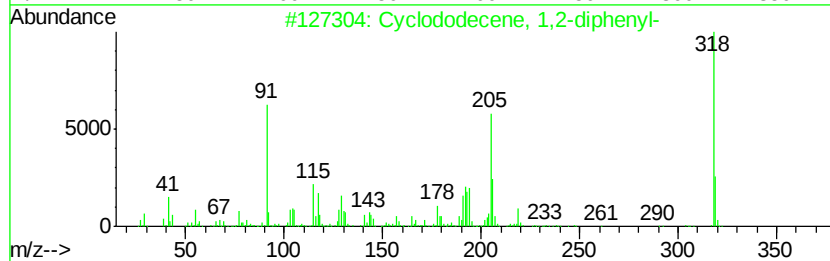
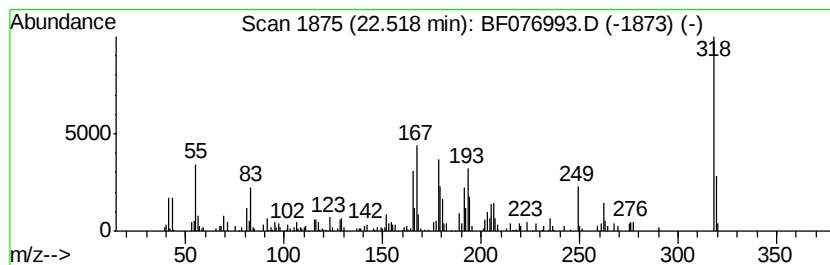
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 unknown22.52 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.52	9.56 ng	148462	Perylene-d12	22.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclododecene, 1,2-diphenyl-	318	C24H30	115977-56-5	47
2		Coronene, 1,2,2a,3,4,4a,5,6,6a,7...	318	C24H30	125379-41-1	38
3		4-Methyl-2-(2-nitro-5-piperidin-...	364	C20H20N4O3	1000294-85-7	38
4		8-Methoxy-11-[3-methylbutyl]-11H...	318	C21H22N2O	155249-60-8	35
5		Vanadium, (.eta.-5-pentamethylcy...	318	C20H27V	1000157-67-2	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076993.D
Acq On : 21 Jan 2015 9:57
Operator : IZ / TP
Sample : G1111-06
Misc : S
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RBDC6-011615

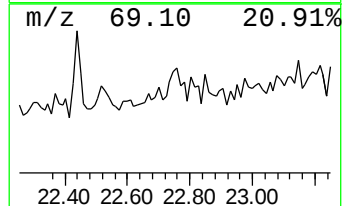
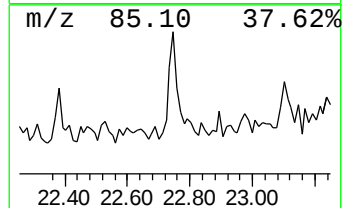
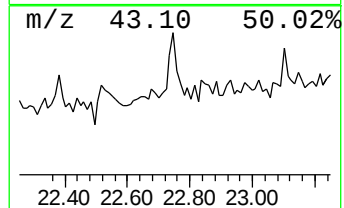
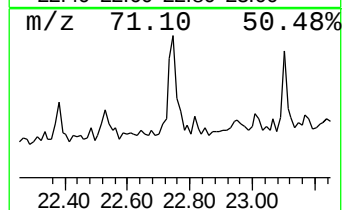
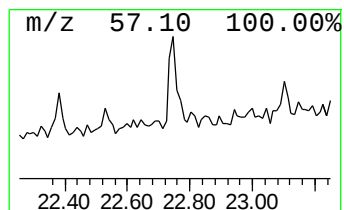
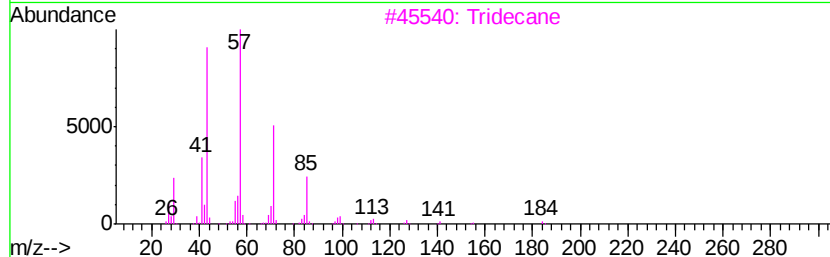
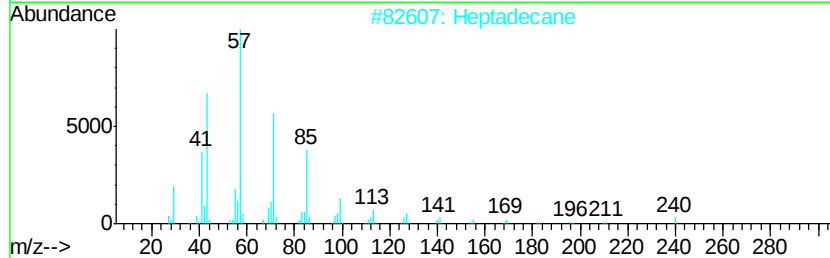
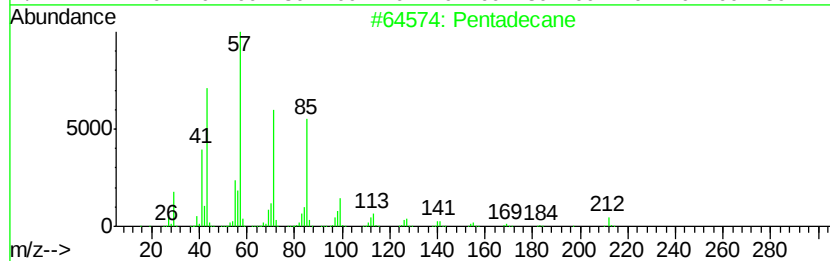
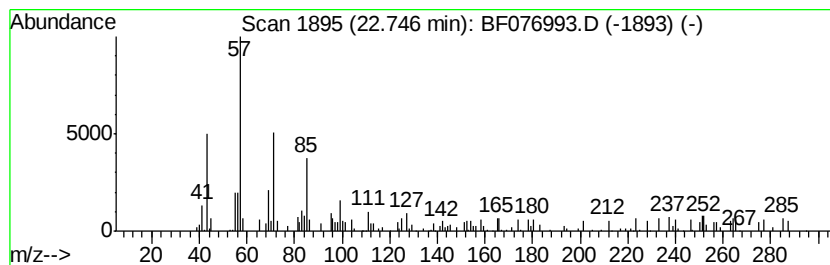
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Pentadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.75	5.61 ng	87068	Perylene-d12	22.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	60
2			Heptadecane	240	C17H36	000629-78-7	49
3			Tridecane	184	C13H28	000629-50-5	47
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	47
5			Tridecane, 2-methyl-	198	C14H30	001560-96-9	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
Data File : BF076993.D
Acq On : 21 Jan 2015 9:57
Operator : IZ / TP
Sample : G1111-06
Misc : S
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RBDC6-011615

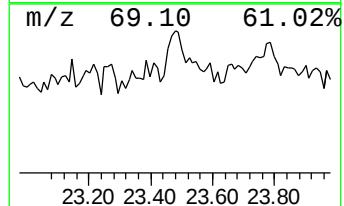
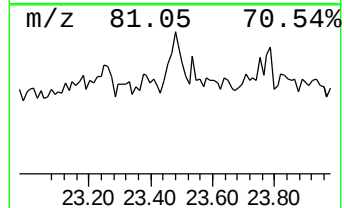
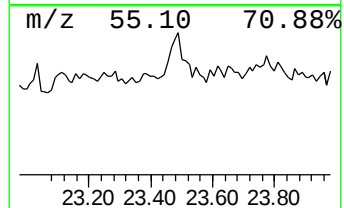
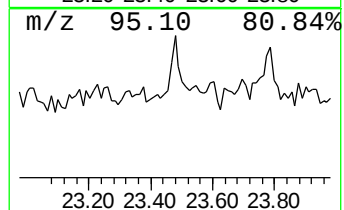
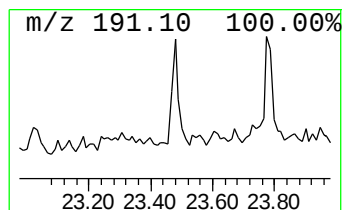
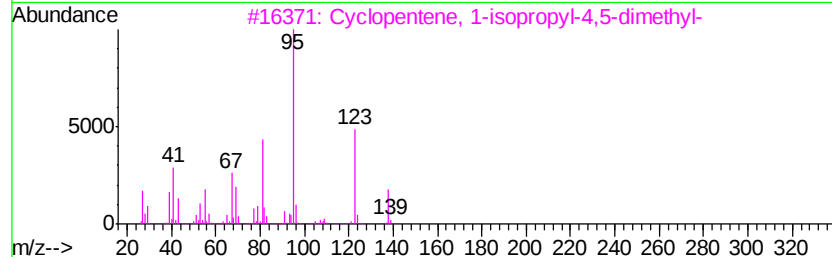
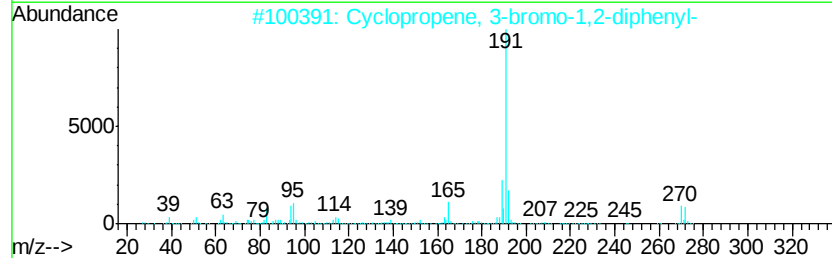
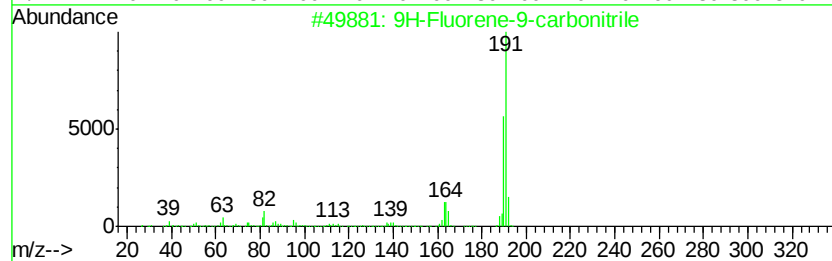
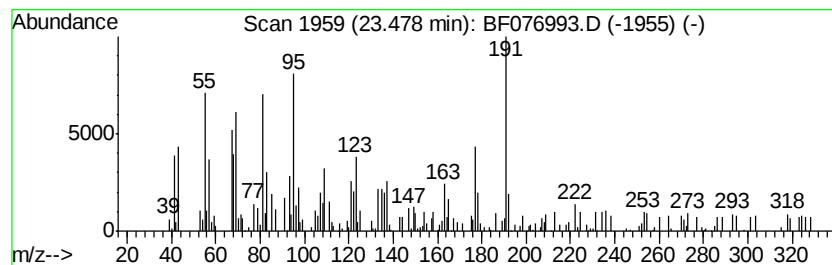
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 unknown23.48 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.48	11.35 ng	176305	Perylene-d12	22.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	43
2		Cyclopropene, 3-bromo-1,2-diphenyl-	270	C15H11Br	029559-17-9	35
3		Cyclopentene, 1-isopropyl-4,5-di...	138	C10H18	007712-74-5	22
4		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	22
5		p-Menth-2-en-9-ol, trans-	154	C10H18O	1000151-16-0	22



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 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBDC6-011615

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.02	15.9	ng	133558	1	7.91	167648	20.0
unknown7.42	7.42	104.6	ng	876417	1	7.91	167648	20.0
Benzophenone	16.59	13.3	ng	190163	4	17.74	286251	20.0
o-Terphenyl	18.37	15.7	ng	224709	4	17.74	286251	20.0
unknown18.54	18.54	9.8	ng	139534	4	17.74	286251	20.0
n-Hexadecanoic acid	18.73	14.6	ng	208244	4	17.74	286251	20.0
Tricyclo[8.2.2.2(...	19.19	20.7	ng	296017	4	17.74	286251	20.0
unknown19.45	19.45	83.3	ng	1192660	4	17.74	286251	20.0
Carbamazepine	19.67	36.3	ng	866463	5	21.24	478011	20.0
unknown22.25	22.25	5.6	ng	86811	6	22.88	310563	20.0
unknown22.52	22.52	9.6	ng	148462	6	22.88	310563	20.0
Pentadecane	22.75	5.6	ng	87068	6	22.88	310563	20.0
unknown23.48	23.48	11.4	ng	176305	6	22.88	310563	20.0