

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
 Data File : BF076989.D
 Acq On : 21 Jan 2015 7:39
 Operator : IZ / TP
 Sample : G1111-02
 Misc : S
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBDC2-011615

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	1.361	23	24	29	rBV2	8417	16569	2.32%	0.200%
2	1.441	29	31	44	rVB	69677	128323	17.93%	1.548%
3	4.024	254	257	266	rBV	47451	109350	15.28%	1.319%
4	4.961	336	339	349	rBV	360227	668940	93.47%	8.070%
5	7.316	541	545	550	rBV	377802	664419	92.84%	8.015%
6	7.407	550	553	558	rVB	392065	715685	100.00%	8.634%
7	7.910	594	597	601	rBV	106339	172738	24.14%	2.084%
8	8.230	621	625	629	rBV	291482	450692	62.97%	5.437%
9	9.202	707	710	717	rBV	261059	405189	56.62%	4.888%
10	10.813	848	851	855	rBV	160835	242779	33.92%	2.929%
11	12.288	977	980	983	rVB	7052	10030	1.40%	0.121%
12	13.465	1080	1083	1087	rBV	348607	541037	75.60%	6.527%
13	14.540	1174	1177	1182	rVB	26144	39407	5.51%	0.475%
14	14.951	1209	1213	1216	rBV	160061	270885	37.85%	3.268%
15	15.397	1249	1252	1256	rBV	8132	17852	2.49%	0.215%
16	16.163	1316	1319	1322	rVB	67657	90963	12.71%	1.097%
17	16.574	1353	1355	1357	rBV	19515	23961	3.35%	0.289%
18	16.631	1357	1360	1363	rVV	329307	420223	58.72%	5.069%
19	17.728	1453	1456	1459	rBV	221526	287883	40.22%	3.473%
20	17.900	1468	1471	1477	rVB2	6015	11594	1.62%	0.140%
21	18.723	1541	1543	1545	rBV	6841	7930	1.11%	0.096%
22	19.420	1601	1604	1607	rVB2	127789	220240	30.77%	2.657%
23	19.694	1625	1628	1631	rVB	33193	35235	4.92%	0.425%
24	19.866	1640	1643	1649	rBV2	2274	7771	1.09%	0.094%
25	19.957	1649	1651	1654	rVV	347990	471098	65.82%	5.683%
26	20.952	1736	1738	1741	rBV3	4456	7625	1.07%	0.092%
27	21.054	1745	1747	1751	rVB3	3036	7163	1.00%	0.086%
28	21.134	1751	1754	1758	rBV3	4724	10526	1.47%	0.127%
29	21.226	1758	1762	1764	rBV	248588	409078	57.16%	4.935%
30	21.260	1764	1765	1767	rVV	22480	22284	3.11%	0.269%
31	21.306	1767	1769	1772	rVV	17065	18703	2.61%	0.226%
32	21.363	1772	1774	1777	rVV	94571	111831	15.63%	1.349%
33	21.615	1792	1796	1798	rBV3	8133	19156	2.68%	0.231%
34	21.660	1798	1800	1803	rVV2	5673	9240	1.29%	0.111%

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35	21.832	1812	1815	1816	rBV3	12889	18453	2.58%	0.223%
36	21.923	1821	1823	1826	rVB	26749	28265	3.95%	0.341%
37	22.003	1827	1830	1836	rVB	74406	96925	13.54%	1.169%
38	22.369	1860	1862	1866	rBV4	11431	20692	2.89%	0.250%
39	22.437	1866	1868	1869	rVV2	6543	12332	1.72%	0.149%
40	22.460	1869	1870	1874	rVV2	21956	43818	6.12%	0.529%
41	22.517	1874	1875	1878	rVV2	16096	23245	3.25%	0.280%
42	22.735	1891	1894	1898	rBV2	83829	140164	19.58%	1.691%
43	22.815	1898	1901	1903	rBV2	16633	26031	3.64%	0.314%
44	22.872	1903	1906	1910	rVV	221917	276031	38.57%	3.330%
45	23.078	1922	1924	1926	rBV2	10260	19278	2.69%	0.233%
46	23.249	1936	1939	1941	rBV	65938	92832	12.97%	1.120%
47	23.386	1948	1951	1953	rBV3	8393	17738	2.48%	0.214%
48	23.432	1953	1955	1957	rBV	33327	63012	8.80%	0.760%
49	23.466	1957	1958	1961	rVB3	39336	50623	7.07%	0.611%
50	23.980	1999	2003	2006	rBV3	25336	65806	9.19%	0.794%
51	24.140	2013	2017	2019	rBV4	13735	39609	5.53%	0.478%
52	24.186	2019	2021	2023	rVV2	12636	21280	2.97%	0.257%
53	24.243	2023	2026	2030	rVB3	22497	53150	7.43%	0.641%
54	24.403	2037	2040	2044	rBV2	82404	172613	24.12%	2.082%
55	24.472	2044	2046	2053	rVB7	22083	62627	8.75%	0.756%
56	24.655	2059	2062	2066	rVB5	25080	50289	7.03%	0.607%
57	24.746	2066	2070	2073	rVB6	11409	26305	3.68%	0.317%
58	24.895	2080	2083	2088	rVB6	26952	73721	10.30%	0.889%
59	25.055	2094	2097	2099	rBV3	12450	28330	3.96%	0.342%
60	25.215	2109	2111	2113	rBV3	4603	8026	1.12%	0.097%
61	25.466	2131	2133	2135	rBV3	4212	7691	1.07%	0.093%
62	25.558	2137	2141	2142	rBV4	6883	15994	2.23%	0.193%
63	25.695	2149	2153	2158	rVB5	23767	59611	8.33%	0.719%
64	26.644	2231	2236	2240	rBV8	5888	20368	2.85%	0.246%
65	27.067	2272	2273	2278	rVB5	3749	8008	1.12%	0.097%

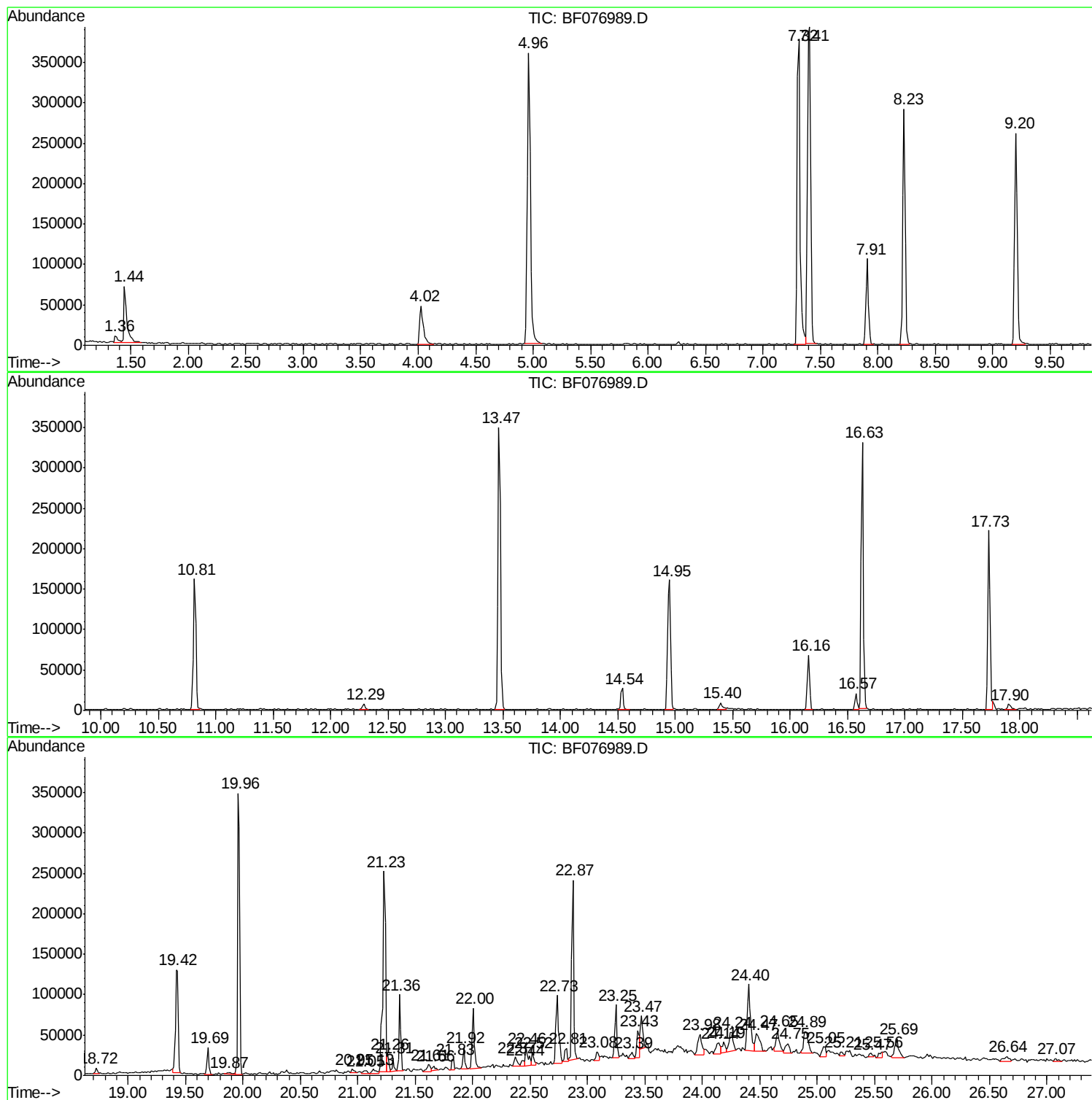
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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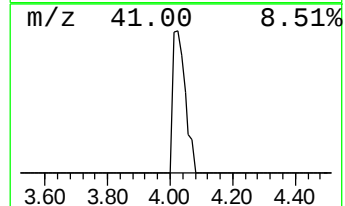
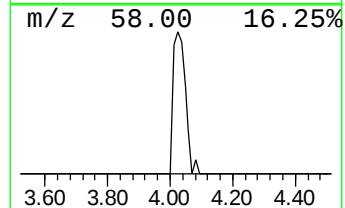
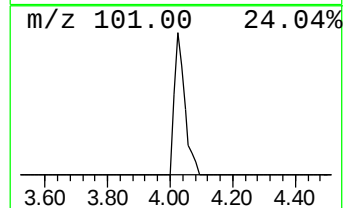
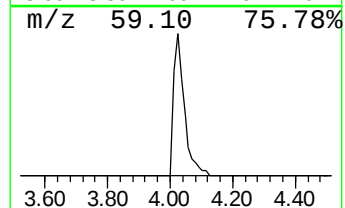
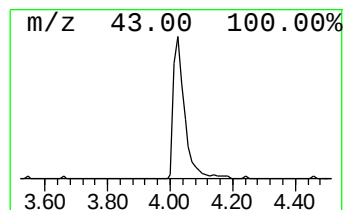
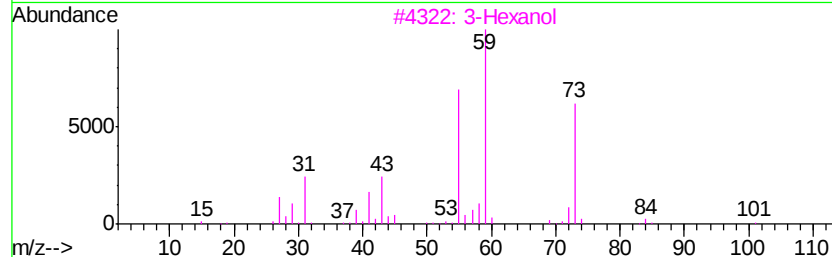
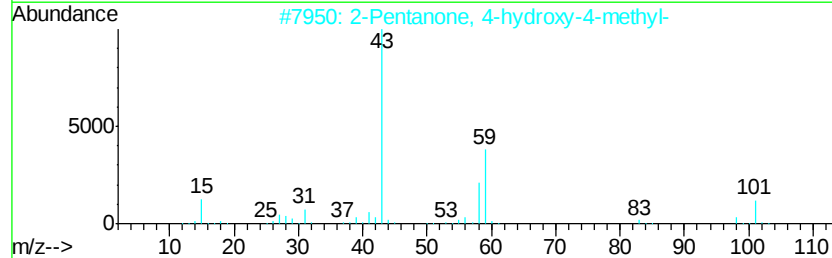
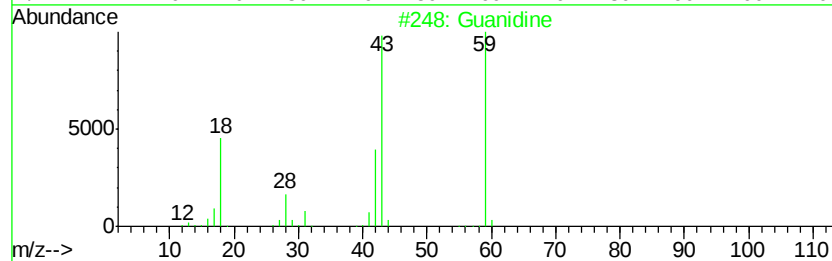
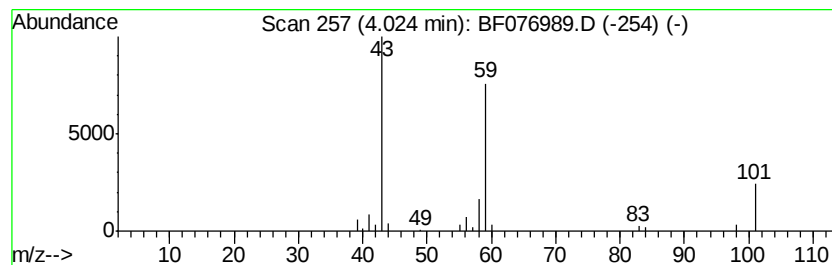
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Guanine Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.02	12.66 ng	109350	1,4-Dichlorobenzene-d4	7.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Guanine		59	CH5N3	000113-00-8	50
2	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	38
3	3-Hexanol		102	C6H14O	000623-37-0	33
4	Butyl aldoxime, 3-methyl-, syn-		101	C5H11NO	005780-40-5	28
5	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	12



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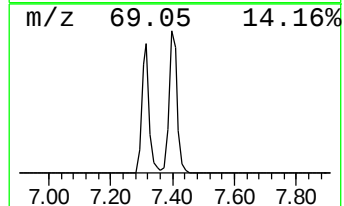
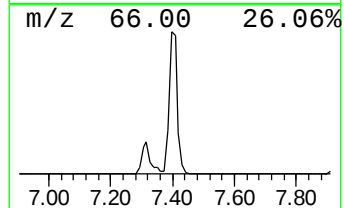
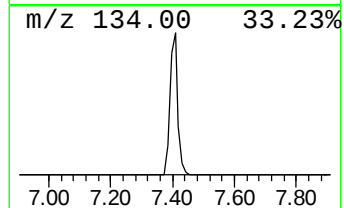
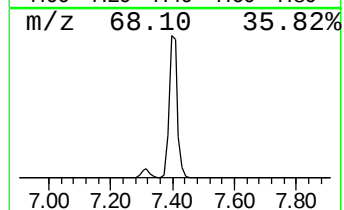
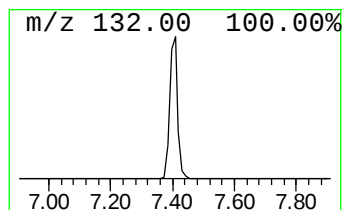
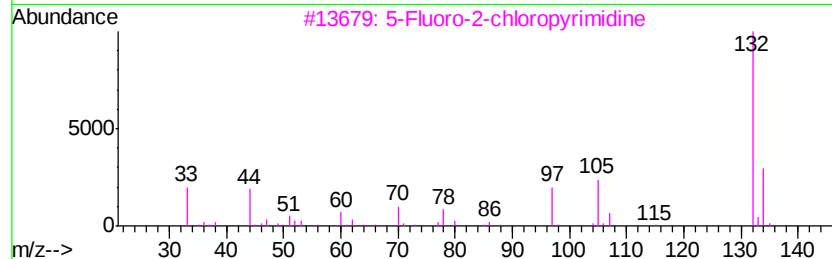
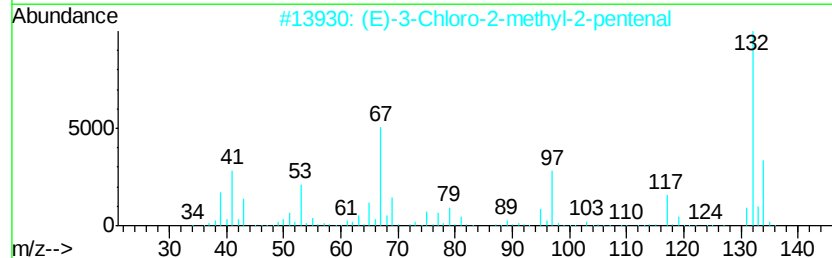
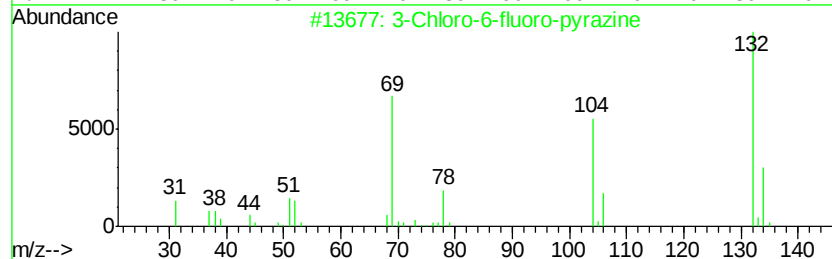
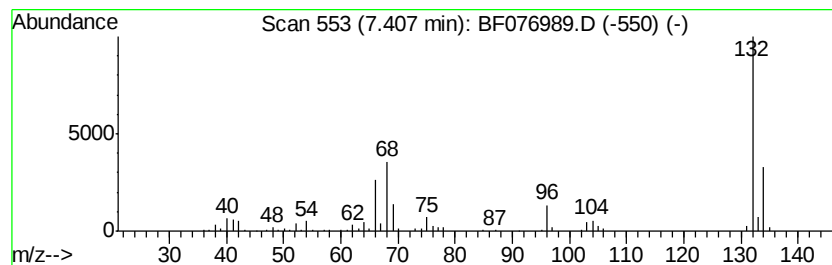
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.41 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.41	82.86 ng	715685	1,4-Dichlorobenzene-d4	7.91

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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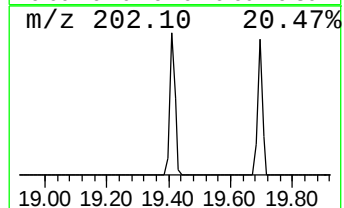
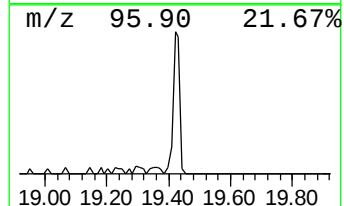
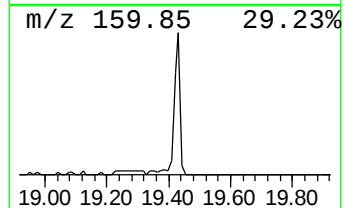
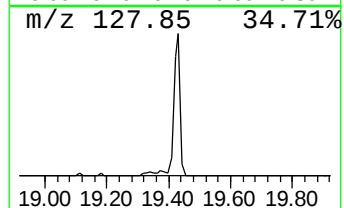
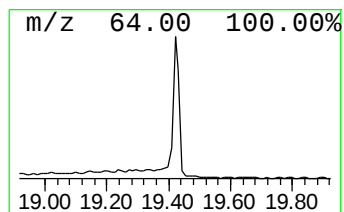
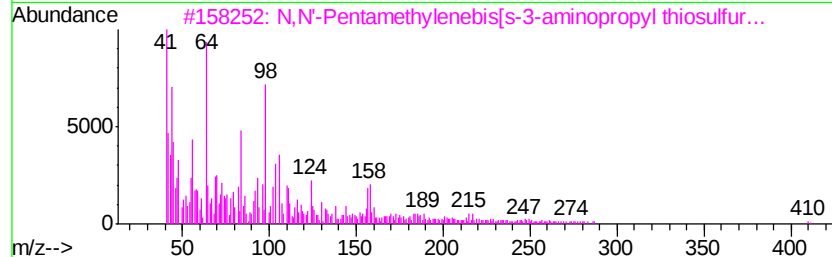
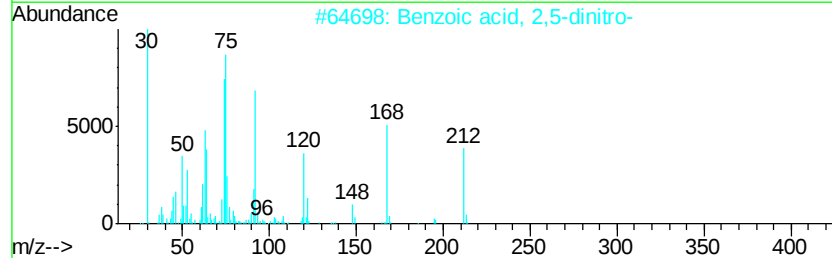
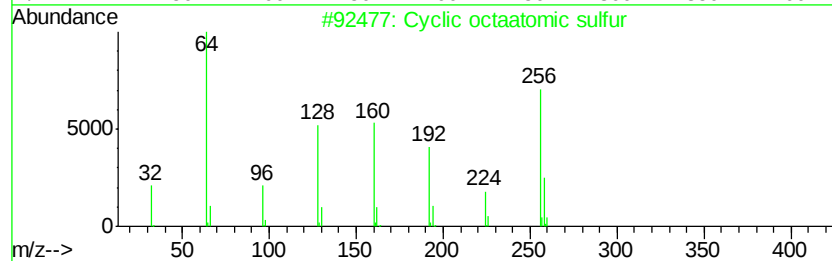
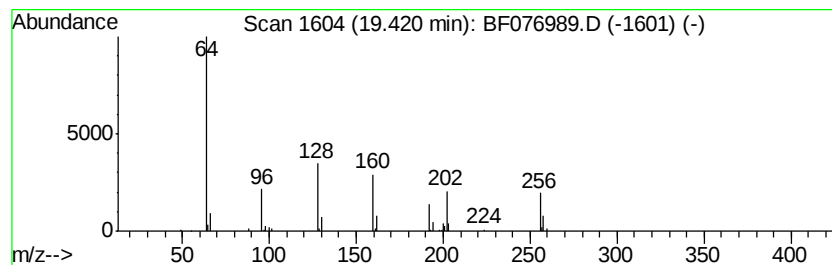
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclic octaatomic sulfur Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.42	15.30 ng	220240	Phenanthrene-d10	17.73

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclic octaatomic sulfur	256	S8	010544-50-0	94
2		Benzoic acid, 2,5-dinitro-	212	C7H4N2O6	000610-28-6	53
3		N,N'-Pentamethylenebis[s-3-amino...	410	C11H26N2O6S4	035871-54-6	38
4		4-Phenylhydrazono-3-methyl-2-pyr...	202	C10H10N4O	1000143-29-5	9
5		2-Benzimidazolyl methane thiosul...	244	C8H8N2O3S2	1000256-39-2	9



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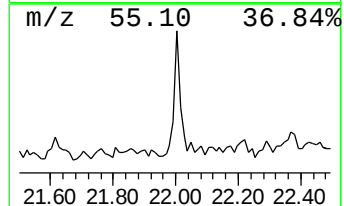
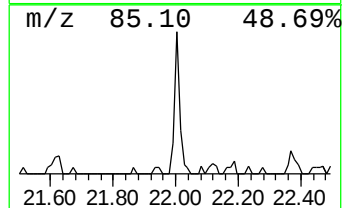
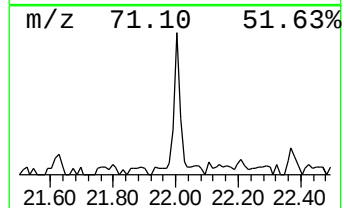
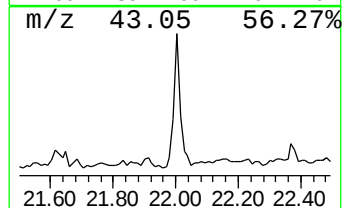
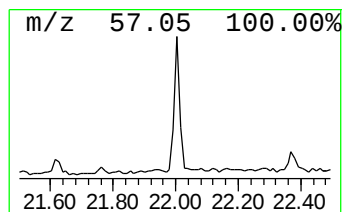
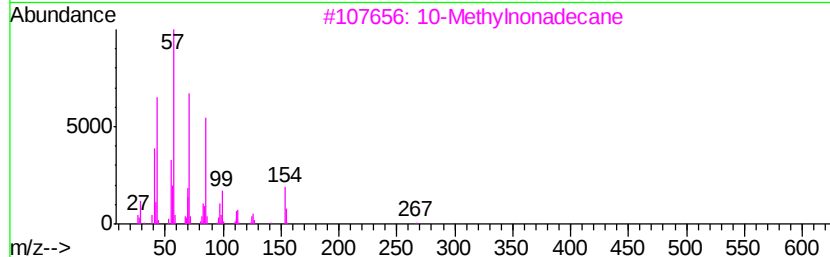
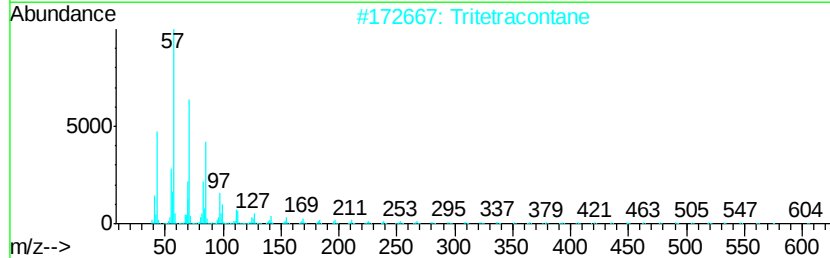
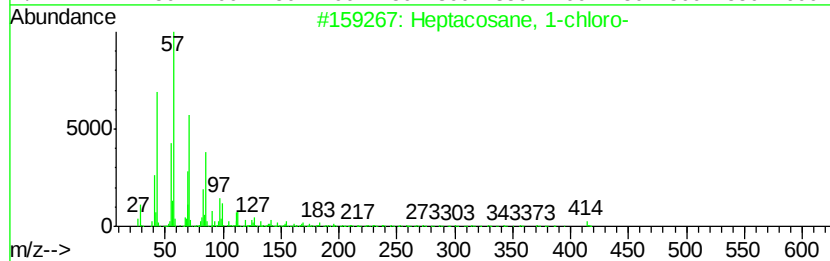
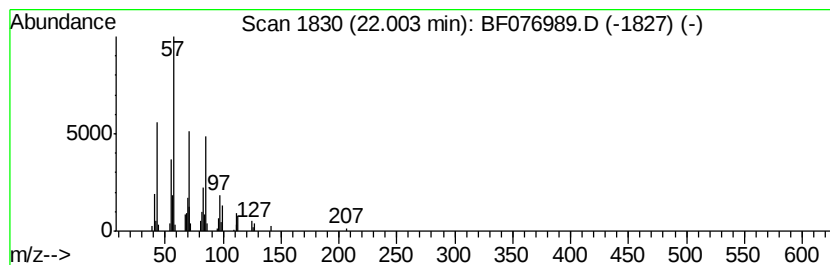
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Heptacosane, 1-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.00	4.74 ng	96925	Chrysene-d12	21.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	90
2			Tritetracontane	605	C43H88	007098-21-7	80
3			10-Methylnonadecane	282	C20H42	056862-62-5	80
4			Nonadecane	268	C19H40	000629-92-5	68
5			Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	64



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RBDC2-011615

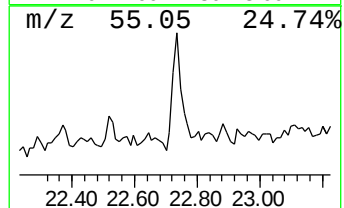
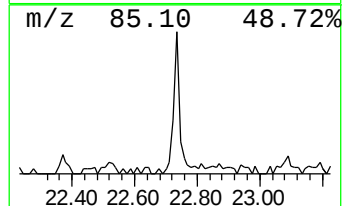
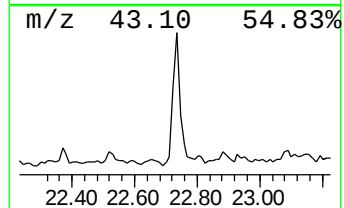
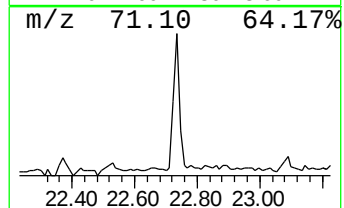
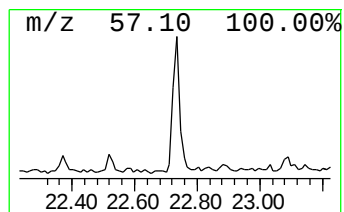
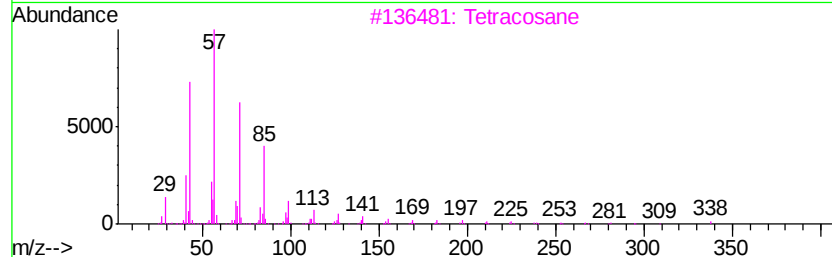
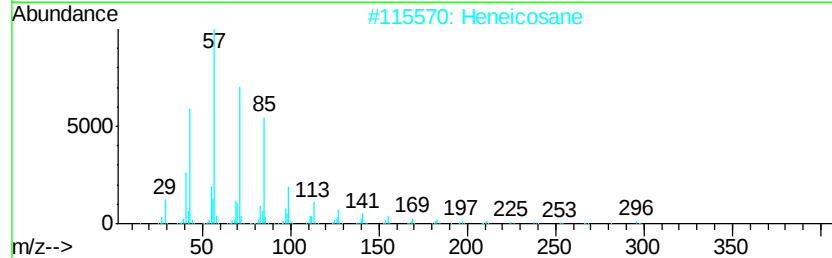
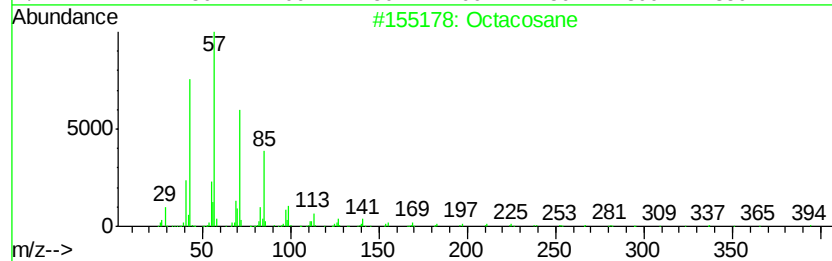
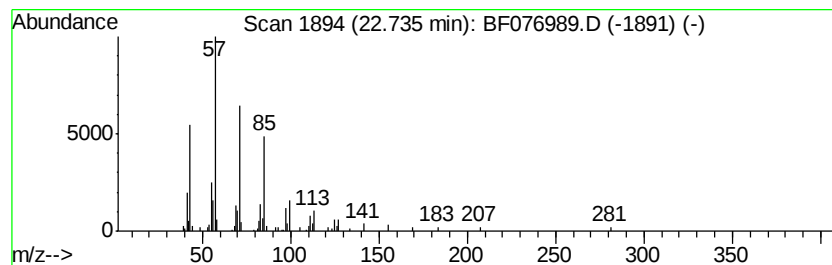
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.73	10.16 ng	140164	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

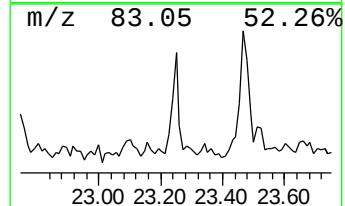
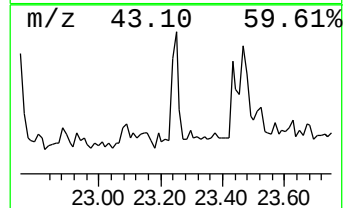
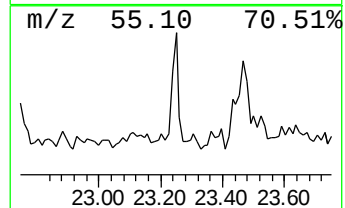
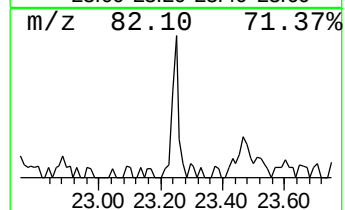
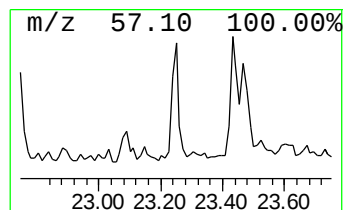
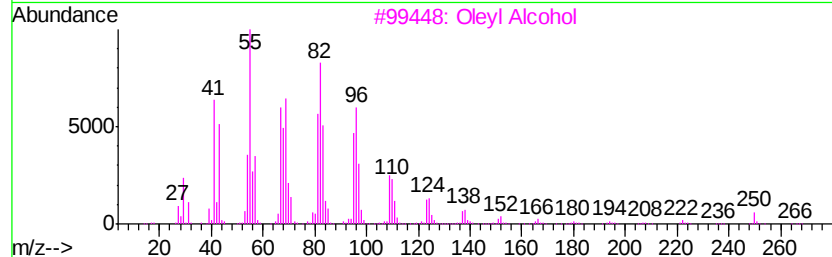
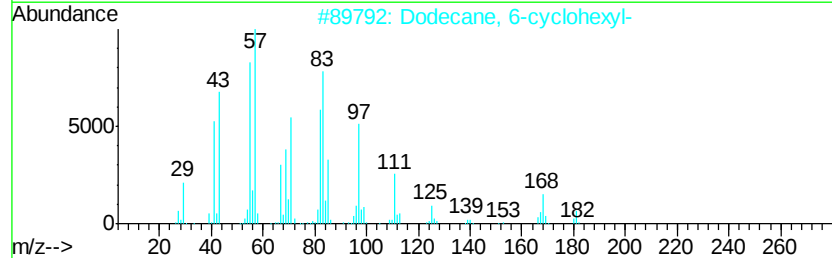
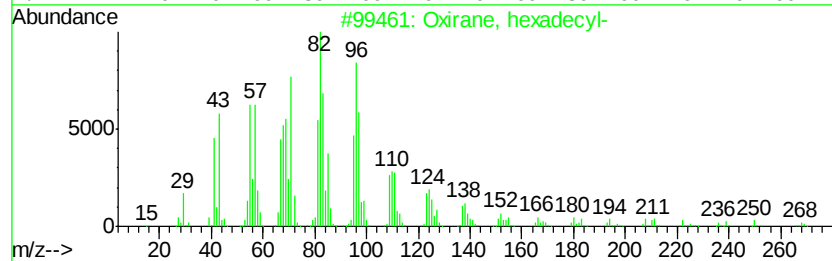
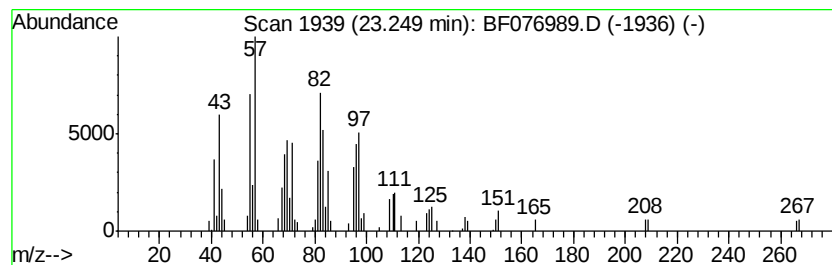
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Oxirane, hexadecyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.25	6.73 ng	92832	Perylene-d12	22.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	64
2	Dodecane, 6-cyclohexyl-	252	C18H36	013151-86-5	58
3	Olevl Alcohol	268	C18H36O	000143-28-2	53
4	Octadecanal	268	C18H36O	000638-66-4	53
5	Ethanol, 2-(9-octadecenyl-oxy)-, ...	312	C20H40O2	005353-25-3	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

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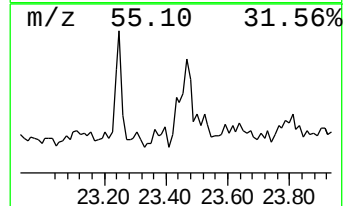
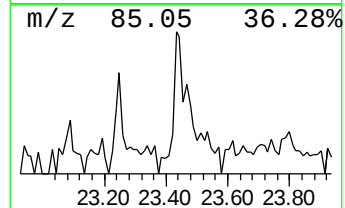
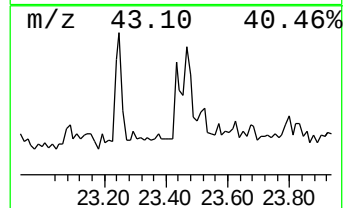
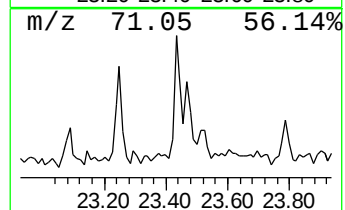
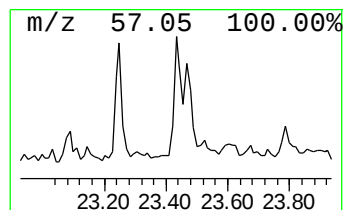
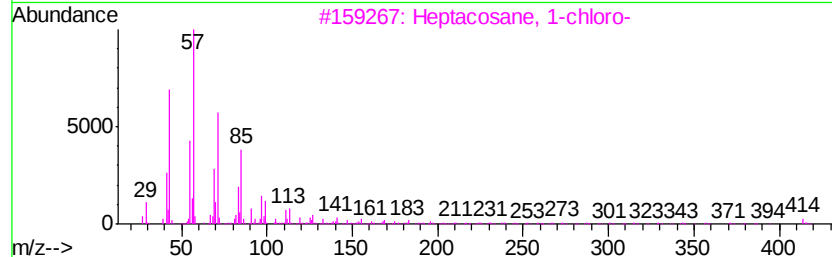
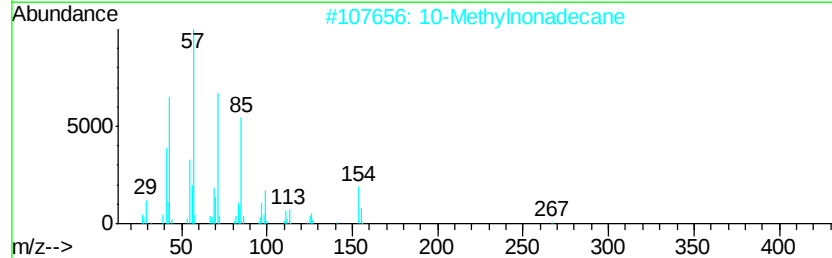
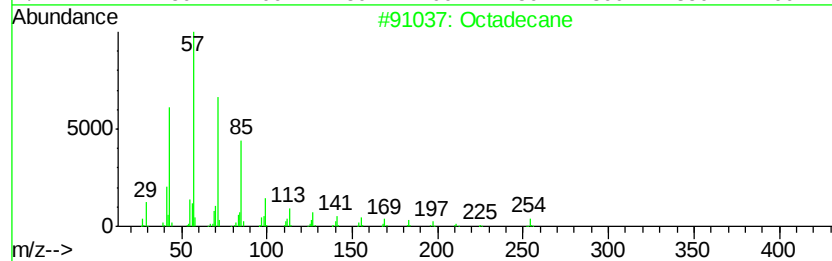
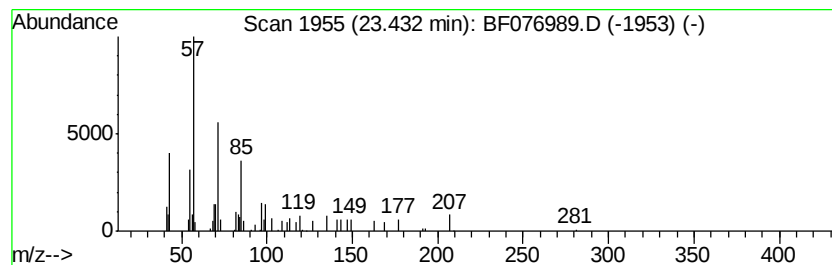
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.43	4.57 ng	63012	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	72
2		10-Methylnonadecane	282	C20H42	056862-62-5	72
3		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	72
4		Docosane	310	C22H46	000629-97-0	64
5		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

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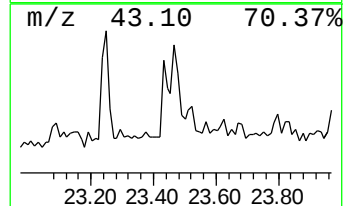
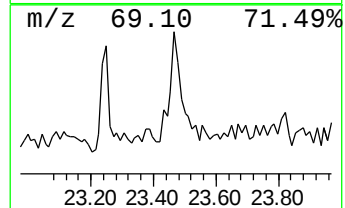
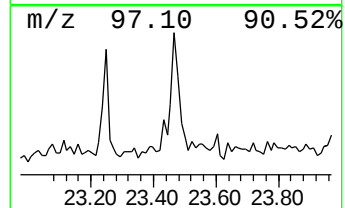
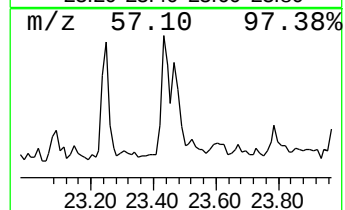
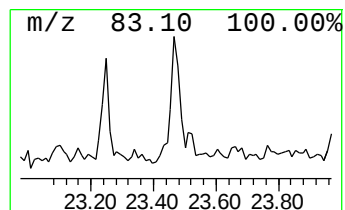
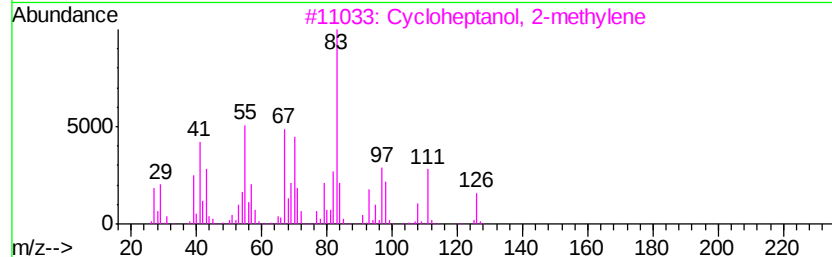
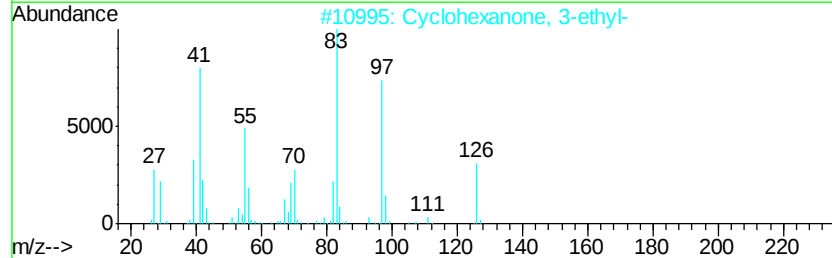
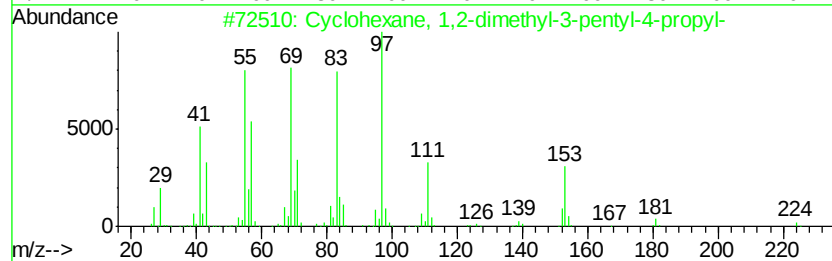
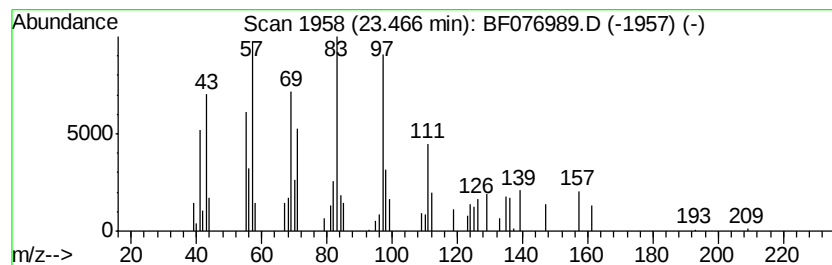
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown23.47 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.47	3.67 ng	50623	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	47
2		Cyclohexanone, 3-ethyl-	126	C8H14O	022461-89-8	43
3		Cycloheptanol, 2-methylene	126	C8H14O	016240-38-3	38
4		Cyclohexanecarboxylic acid, 4-oc...	240	C15H28O2	1000279-52-2	35
5		Cyclohexane, 1,2-dimethyl-	112	C8H16	000583-57-3	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

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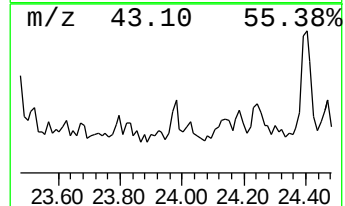
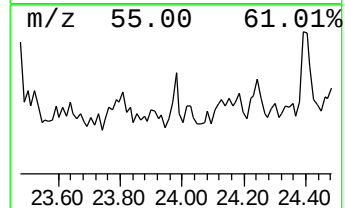
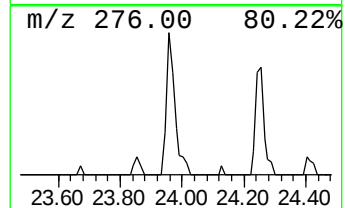
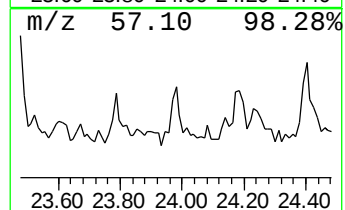
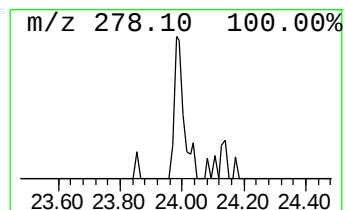
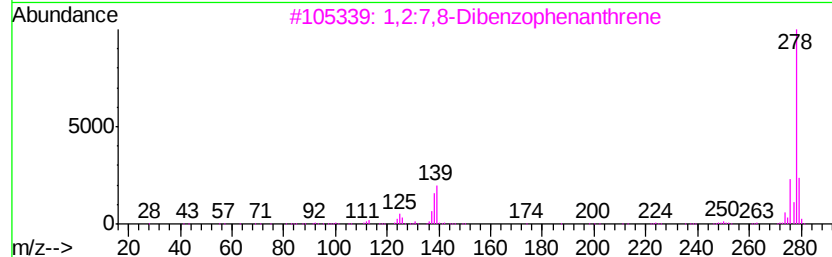
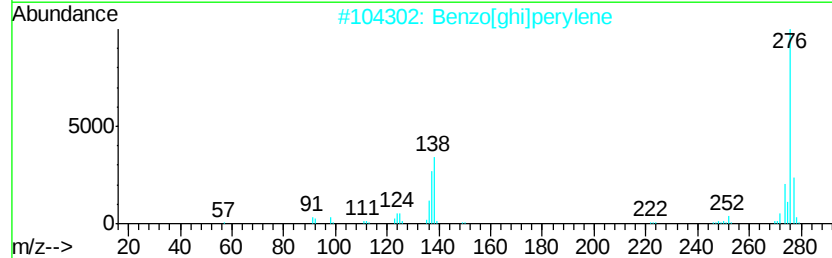
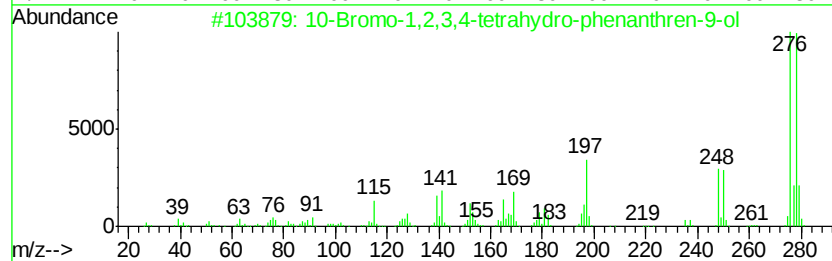
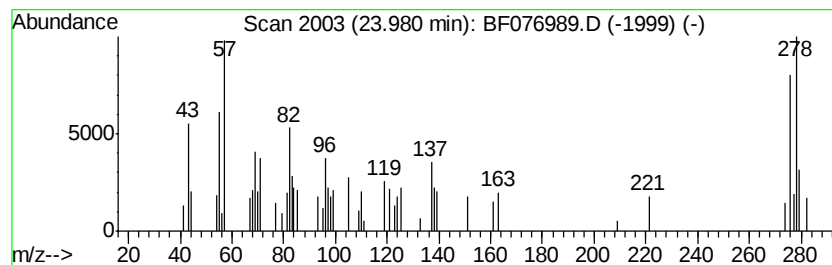
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown23.98 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.98	4.77 ng	65806	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Bromo-1,2,3,4-tetrahydro-phen...	276	C14H13BrO	1000190-54-9	40
2		Benzo[ghi]perylene	276	C22H12	000191-24-2	38
3		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	38
4		Benzo[bl]triphenylene	278	C22H14	000215-58-7	38
5		1,3-Diamino-7,8-dichlorobenzo[f]...	278	C12H8Cl2N4	037521-57-6	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

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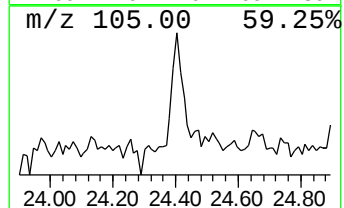
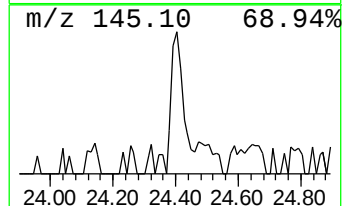
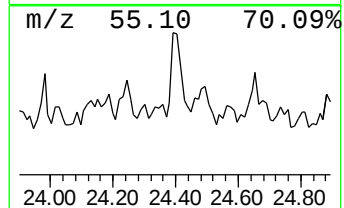
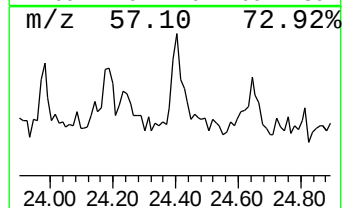
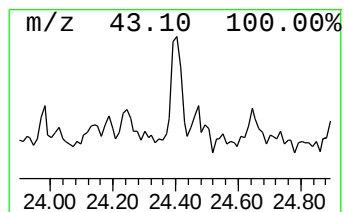
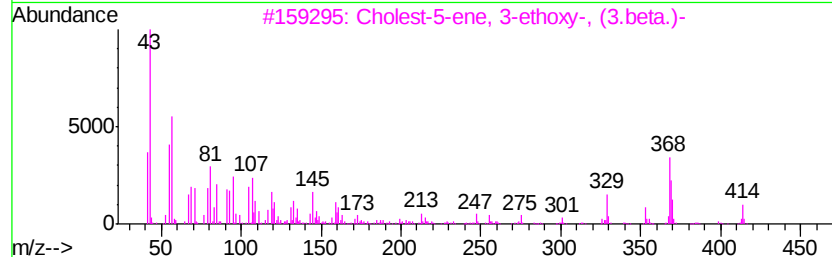
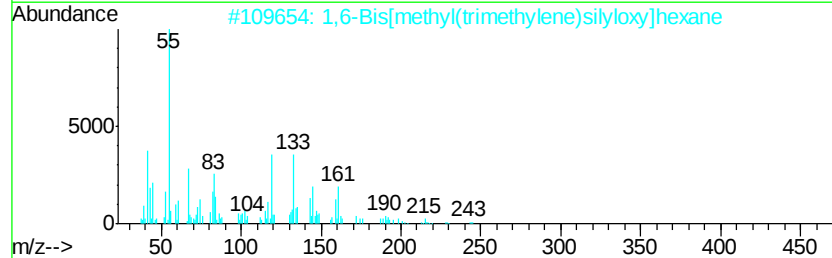
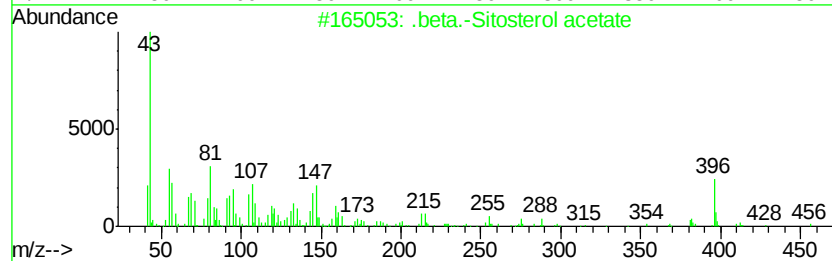
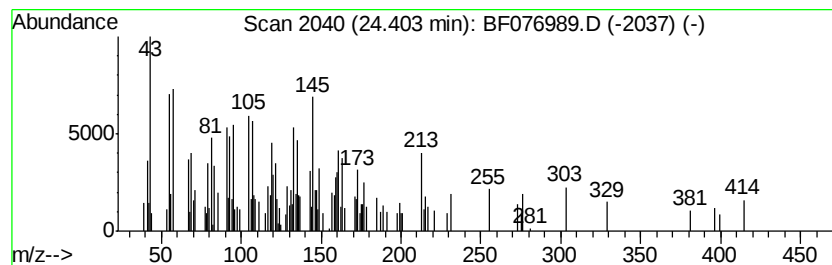
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown24.40 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.40	12.51 ng	172613	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	32
2		1,6-Bis[methyl(trimethylene)silyl]hexane	286	C14H30O2Si2	1000217-00-5	27
3		Cholest-5-ene, 3-ethoxy-, (3.beta.)-	414	C29H50O	000986-19-6	25
4		Cyclopropanecarboxylic acid, 3-(...)	372	C22H28O5	000121-29-9	22
5		Bicyclo[2.2.1]heptane, 2-cyclopr...	176	C13H20	1000159-45-7	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
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Acq On : 21 Jan 2015 7:39
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Misc : S
ALS Vial : 27 Sample Multiplier: 1

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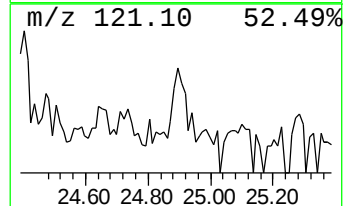
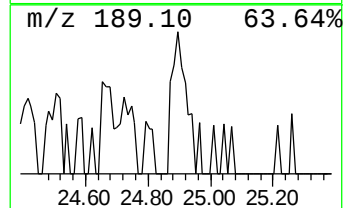
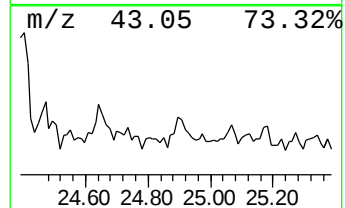
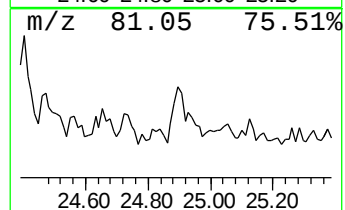
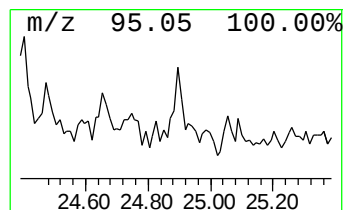
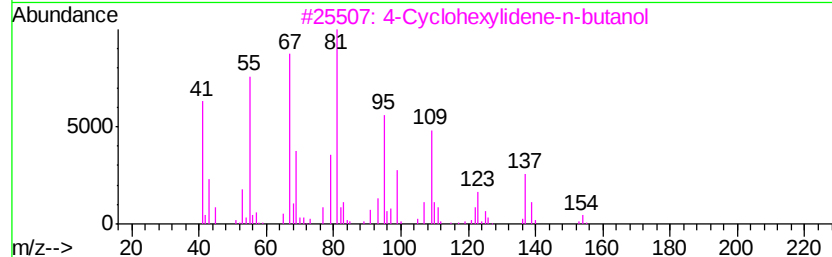
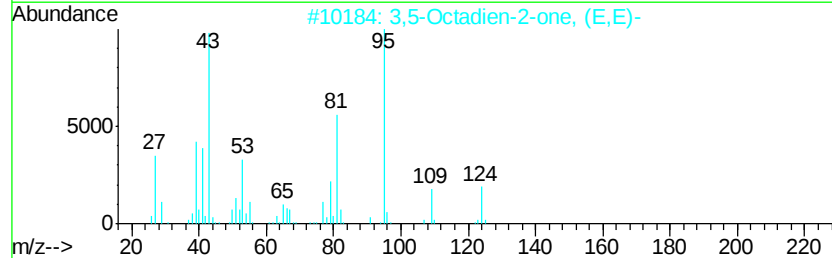
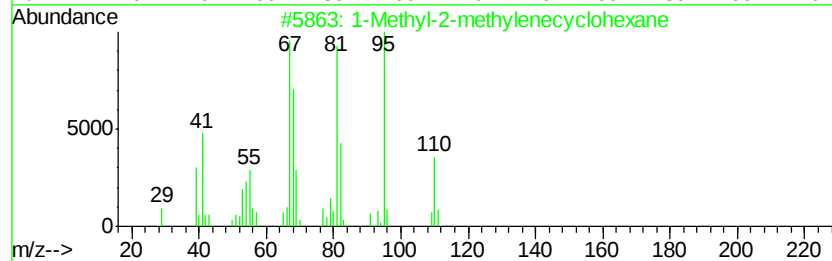
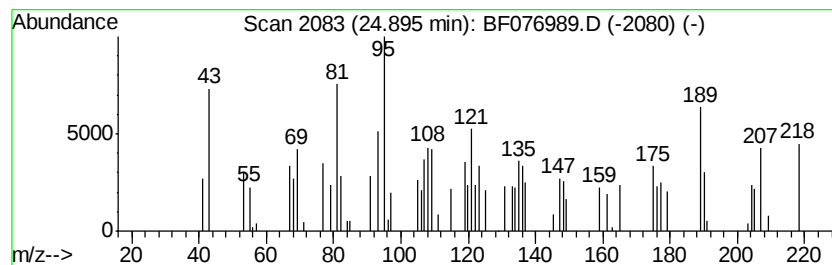
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown24.89 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.89	5.34 ng	73721	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyl-2-methylenecyclohexane	110	C8H14	002808-75-5	27
2		3,5-Octadien-2-one, (E,E)-	124	C8H12O	030086-02-3	27
3		4-Cyclohexylidene-n-butanol	154	C10H18O	004441-58-1	22
4		Thuiopsene-(I2)	204	C15H24	1000152-20-4	15
5		Isobornyl propionate	210	C13H22O2	002756-56-1	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RBDC2-011615

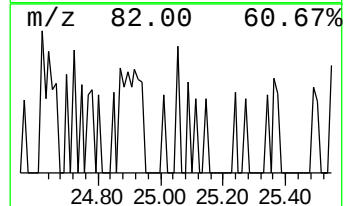
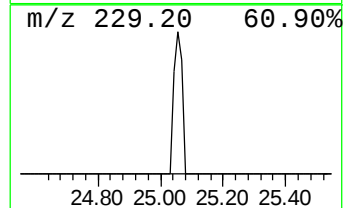
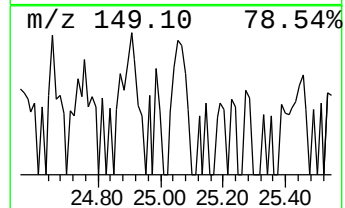
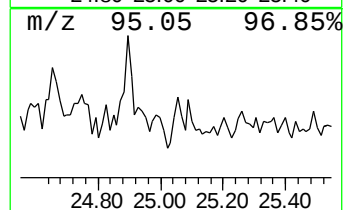
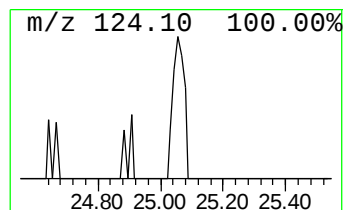
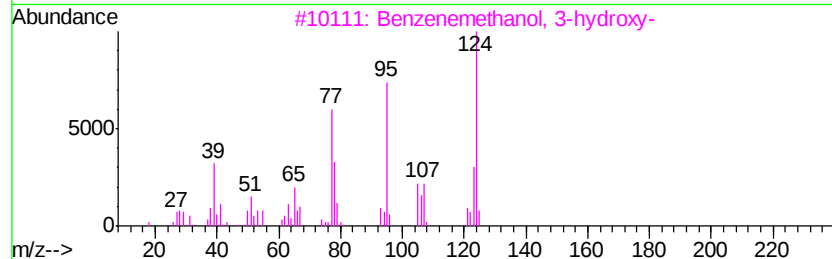
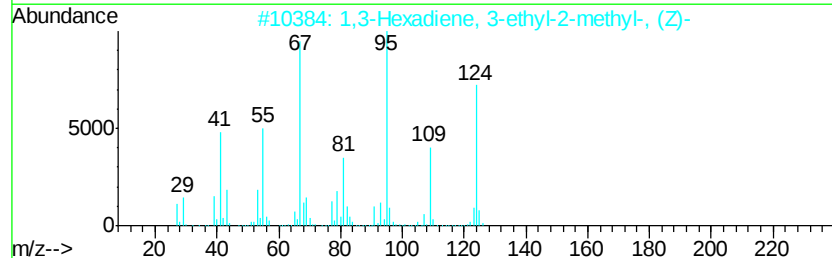
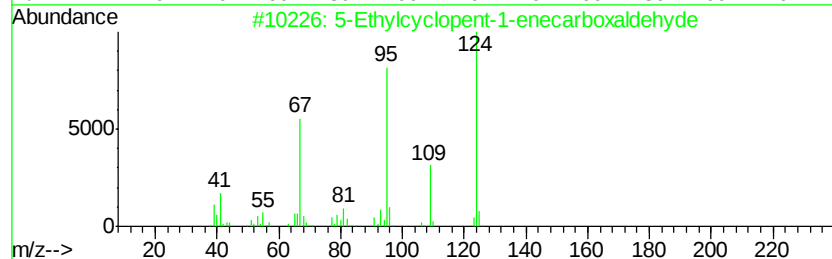
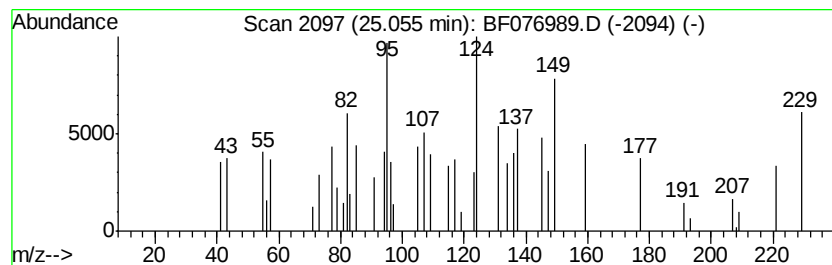
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown25.05 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.05	2.05 ng	28330	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Ethylcyclopent-1-enecarboxalde...	124	C8H12O	036431-60-4	25
2		1,3-Hexadiene, 3-ethyl-2-methyl-...	124	C9H16	074752-97-9	22
3		Benzenemethanol, 3-hydroxy-	124	C7H8O2	000620-24-6	18
4		2(1H)-Pyridinone, 1-propyl-	137	C8H11NO	019006-63-4	18
5		Pyrazine, 2-methoxy-3-methyl-	124	C6H8N2O	002847-30-5	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
Data File : BF076989.D
Acq On : 21 Jan 2015 7:39
Operator : IZ / TP
Sample : G1111-02
Misc : S
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RBDC2-011615

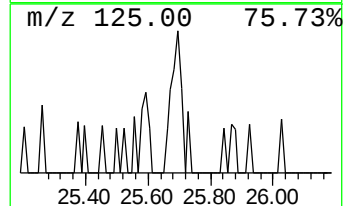
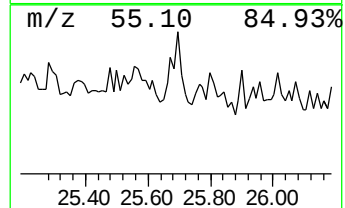
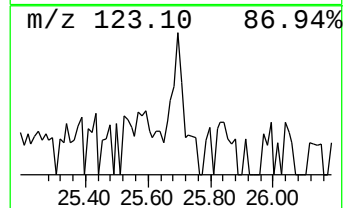
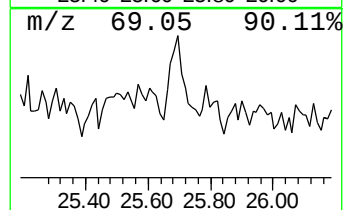
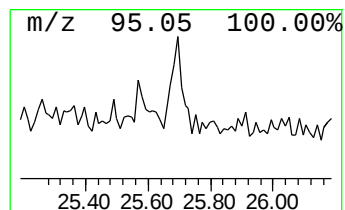
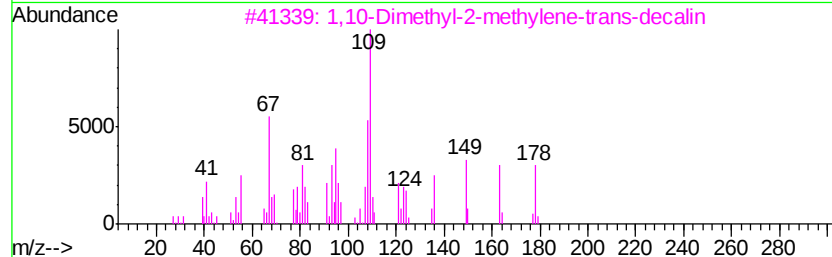
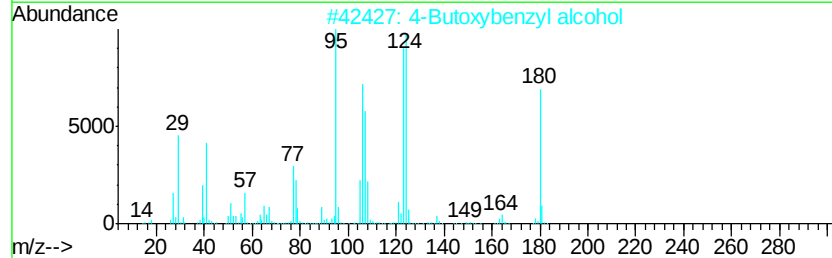
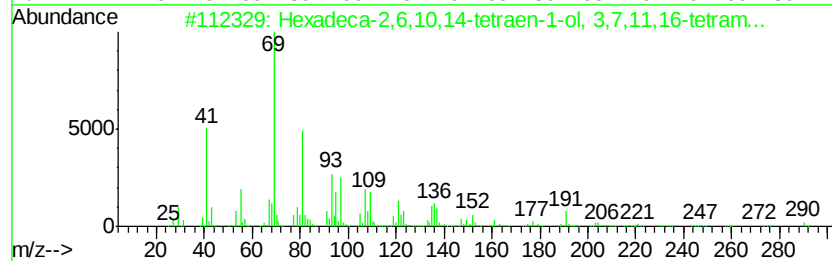
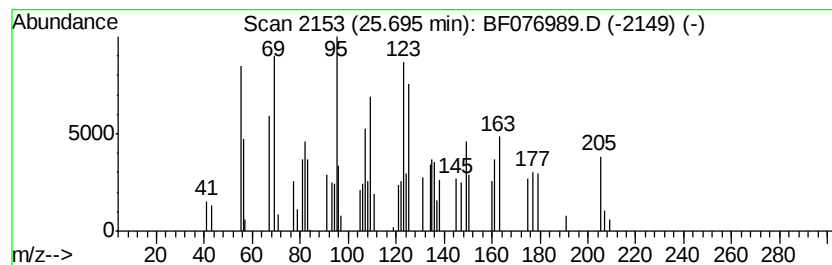
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown25.69 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.69	4.32 ng	59611	Perylene-d12	22.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	22
2		4-Butoxybenzyl alcohol	180	C11H16O2	006214-45-5	22
3		1,10-Dimethyl-2-methylene-trans-...	178	C13H22	1000103-97-8	14
4		6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	12
5		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012015\
 Data File : BF076989.D
 Acq On : 21 Jan 2015 7:39
 Operator : IZ / TP
 Sample : G1111-02
 Misc : S
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBDC2-011615

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
Guanidine	4.02	12.7	ng	109350	1	7.91	172738	20.0
unknown7.41	7.41	82.9	ng	715685	1	7.91	172738	20.0
Cyclic octaatomic...	19.42	15.3	ng	220240	4	17.73	287883	20.0
Heptacosane, 1-ch...	22.00	4.7	ng	96925	5	21.23	409078	20.0
Octacosane	22.73	10.2	ng	140164	6	22.87	276031	20.0
Oxirane, hexadecyl-	23.25	6.7	ng	92832	6	22.87	276031	20.0
Octadecane	23.43	4.6	ng	63012	6	22.87	276031	20.0
unknown23.47	23.47	3.7	ng	50623	6	22.87	276031	20.0
unknown23.98	23.98	4.8	ng	65806	6	22.87	276031	20.0
unknown24.40	24.40	12.5	ng	172613	6	22.87	276031	20.0
unknown24.47	24.47	4.5	ng	62627	6	22.87	276031	20.0
unknown24.89	24.89	5.3	ng	73721	6	22.87	276031	20.0
unknown25.05	25.05	2.0	ng	28330	6	22.87	276031	20.0
unknown25.69	25.69	4.3	ng	59611	6	22.87	276031	20.0