

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110918\
 Data File : BN003486.D
 Acq On : 09 Nov 2018 23:27
 Operator : MJ/SJ
 Sample : J5860-21
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EP2-SB-110(14-16)

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.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	3.028	3	7	19	rVB	93419	116372	1.75%	0.286%
2	3.564	95	98	116	rVB	48320	68741	1.03%	0.169%
3	4.963	331	336	343	rBV	79216	108983	1.63%	0.268%
4	5.416	406	413	426	rVB	2032869	2895111	43.43%	7.127%
5	7.010	675	684	700	rBV	1992396	3446508	51.70%	8.484%
6	7.381	739	747	759	rBV	2122077	3390301	50.85%	8.346%
7	7.846	820	826	832	rBV	358127	563779	8.46%	1.388%
8	8.169	873	881	889	rVB	1629069	2563766	38.46%	6.311%
9	9.016	1016	1025	1033	rBV	1229670	2056133	30.84%	5.062%
10	10.646	1295	1302	1308	rBV	509304	835545	12.53%	2.057%
11	11.634	1465	1470	1475	rBV	48206	70488	1.06%	0.174%
12	12.522	1617	1621	1627	rVB2	59586	89098	1.34%	0.219%
13	12.993	1697	1701	1706	rVB	92222	120702	1.81%	0.297%
14	13.104	1713	1720	1727	rBV	3538223	5253470	78.80%	12.933%
15	13.287	1741	1751	1760	rBV2	105062	221105	3.32%	0.544%
16	13.934	1856	1861	1867	rVV	586284	838041	12.57%	2.063%
17	14.357	1928	1933	1939	rBV	171102	230231	3.45%	0.567%
18	14.487	1949	1955	1963	rVB2	786436	1198735	17.98%	2.951%
19	14.975	2031	2038	2047	rBV7	43649	112810	1.69%	0.278%
20	15.310	2090	2095	2100	rVB	190669	229773	3.45%	0.566%
21	15.710	2156	2163	2168	rBV	77996	148384	2.23%	0.365%
22	15.975	2202	2208	2217	rVB	2959231	4171805	62.58%	10.270%
23	16.169	2236	2241	2243	rBV	196481	262751	3.94%	0.647%
24	16.963	2371	2376	2380	rBV	145676	179365	2.69%	0.442%
25	17.010	2380	2384	2389	rVB2	65901	92843	1.39%	0.229%
26	17.228	2414	2421	2427	rBV	993885	1392708	20.89%	3.428%
27	17.698	2497	2501	2506	rVB	137520	152230	2.28%	0.375%
28	18.098	2564	2569	2575	rBV	184183	256642	3.85%	0.632%
29	18.392	2614	2619	2623	rBV	94883	115473	1.73%	0.284%
30	19.022	2722	2726	2732	rVB	75792	83642	1.25%	0.206%
31	19.375	2782	2786	2797	rBV	48148	80084	1.20%	0.197%
32	19.845	2860	2866	2872	rBV	5301291	6666857	100.00%	16.412%
33	21.098	3074	3079	3085	rBV	120771	166553	2.50%	0.410%
34	21.404	3126	3131	3137	rBV	1048130	1268954	19.03%	3.124%

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EP2-SB-110(14-16)

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	23.721	3517	3525	3535	rVB	623732	1173706	17.61%	2.889%
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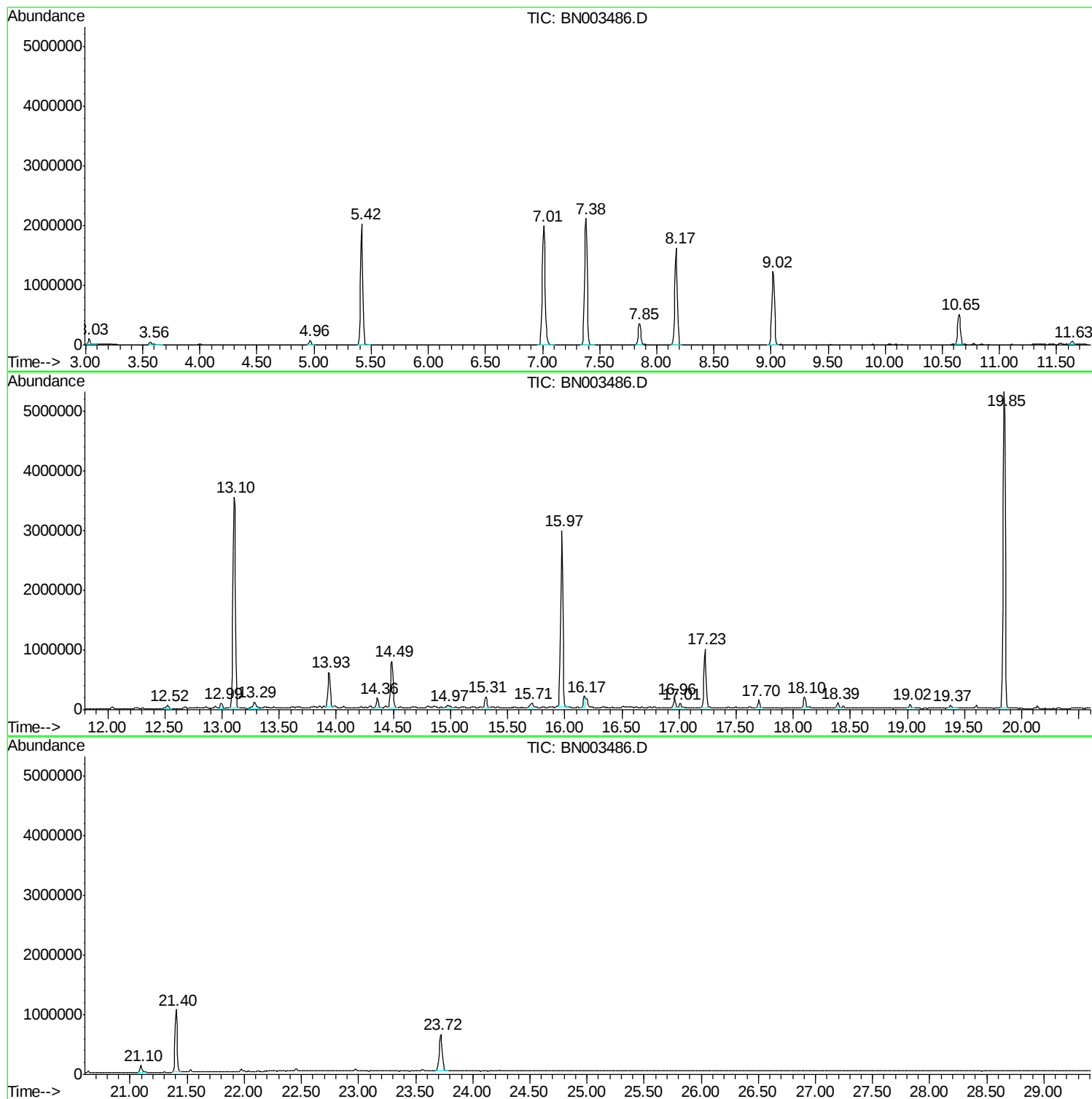
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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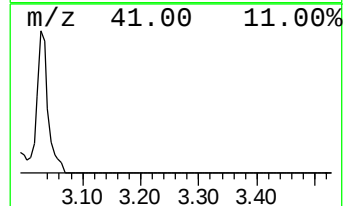
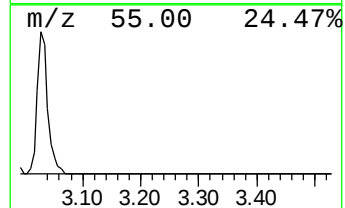
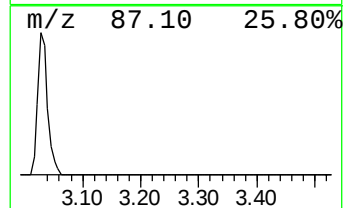
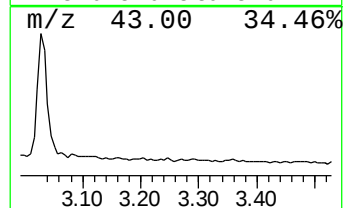
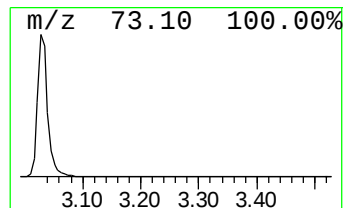
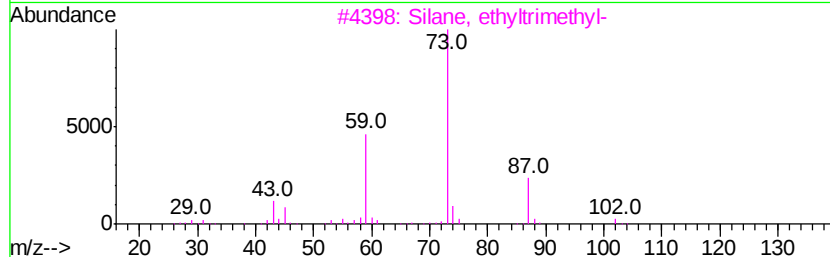
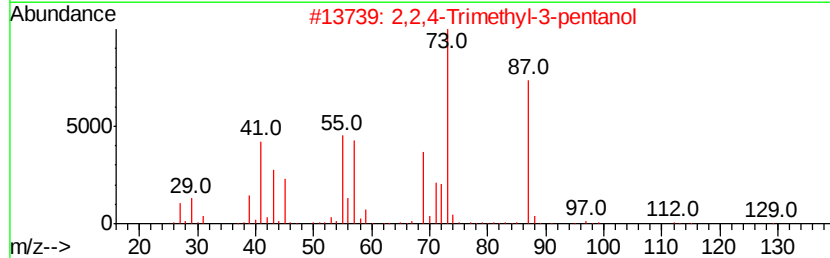
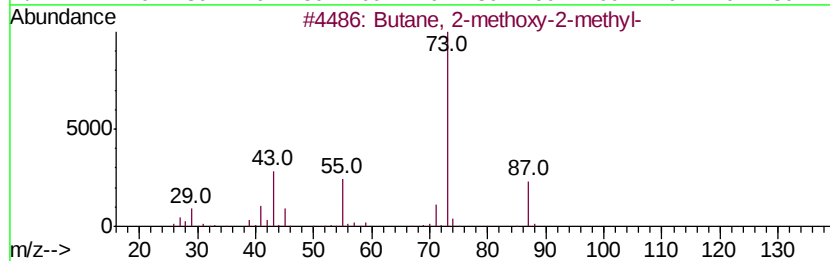
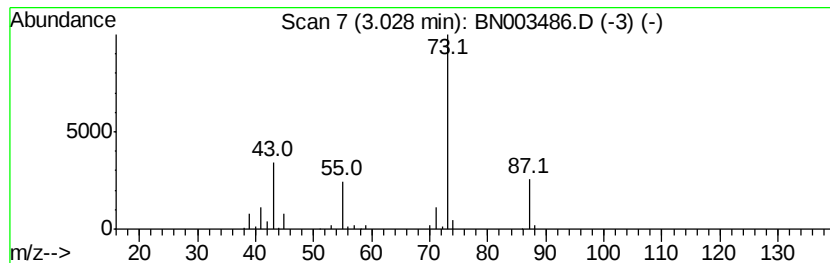
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.03	4.13 ng	116372	1,4-Dichlorobenzene-d4	7.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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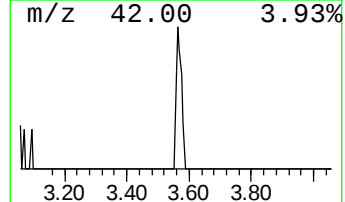
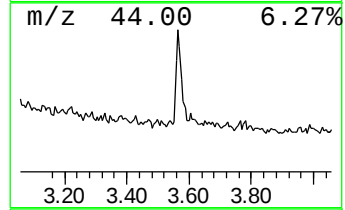
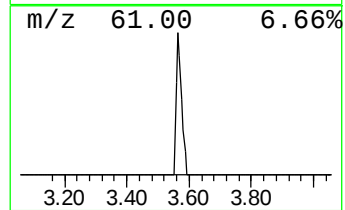
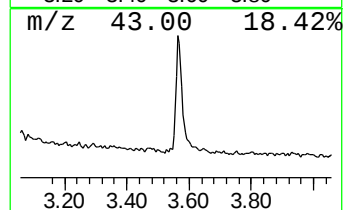
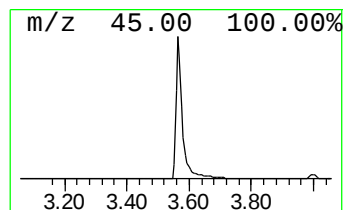
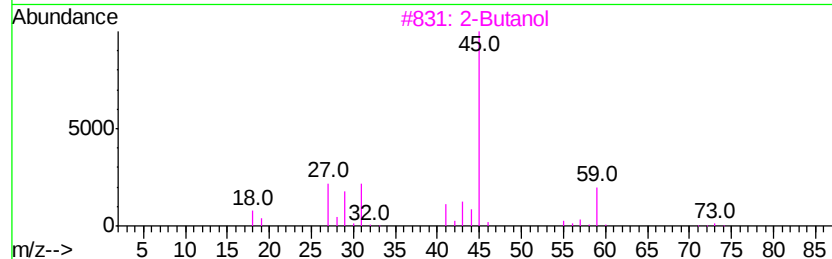
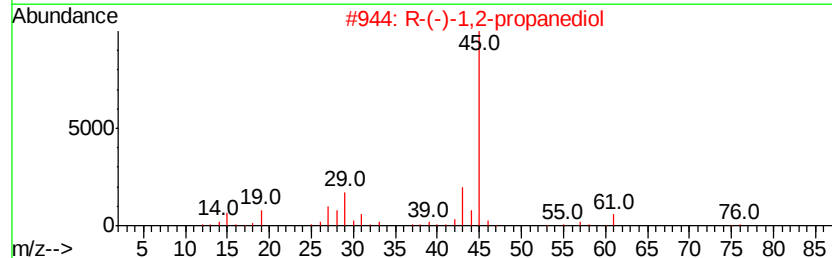
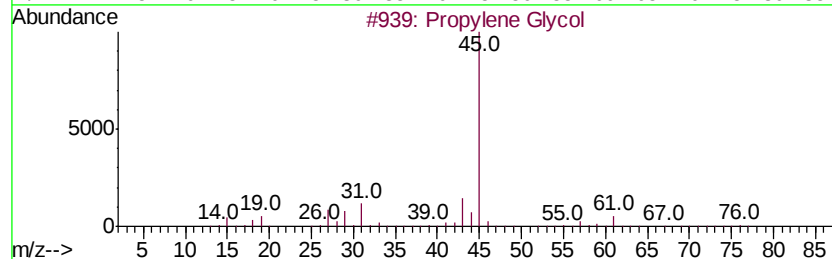
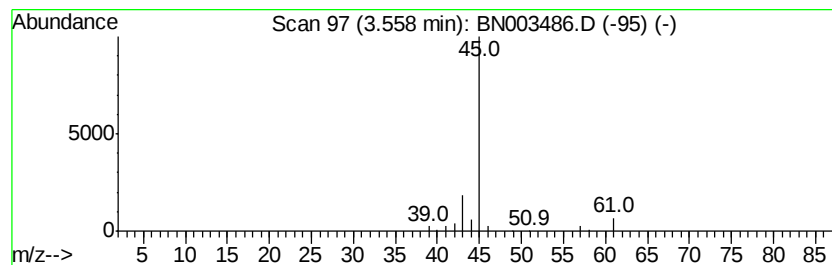
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown3.56 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.56	2.44 ng	68741	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propylene Glycol	76	C3H8O2	000057-55-6	43
2		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
3		2-Butanol	74	C4H10O	000078-92-2	9
4		Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5		Methyltartronic acid	134	C4H6O5	000595-98-2	9



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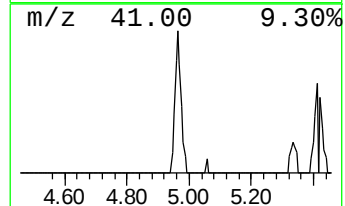
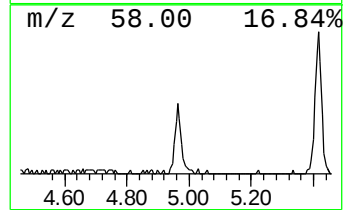
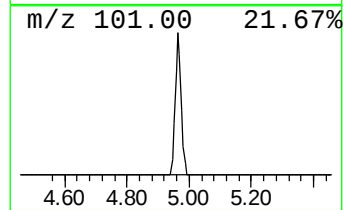
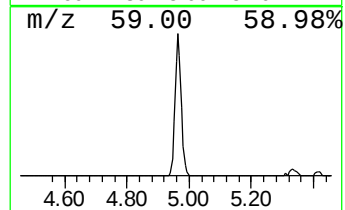
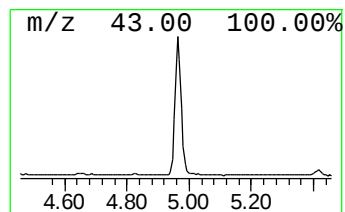
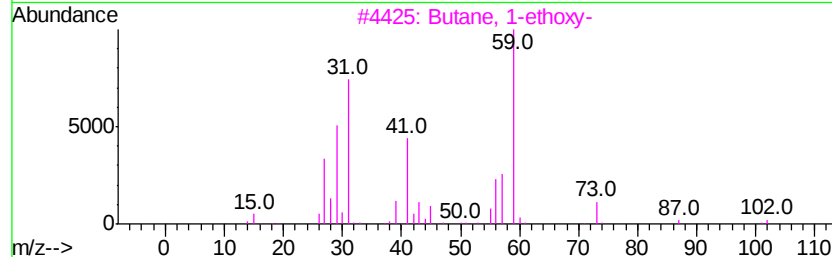
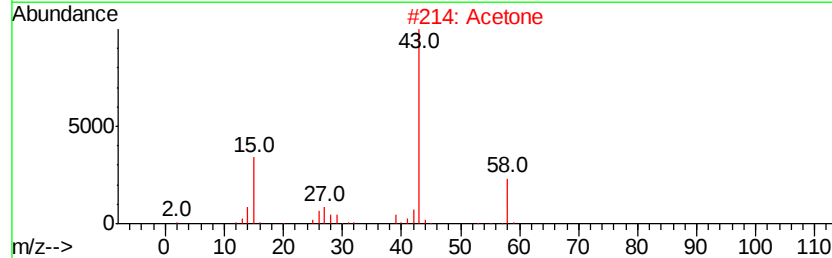
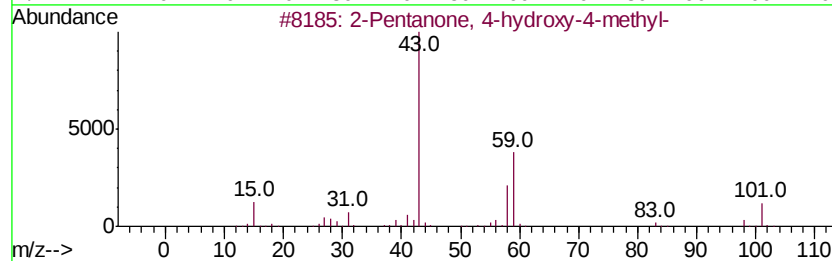
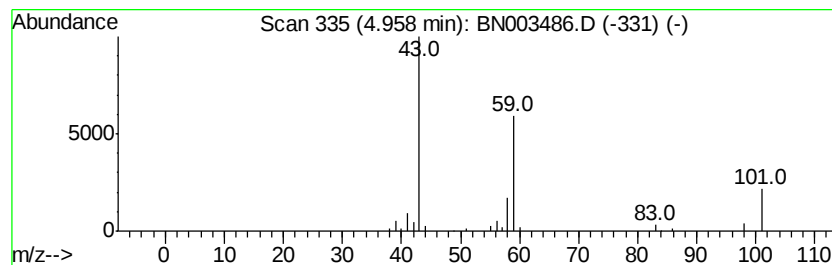
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	3.87 ng	108983	1,4-Dichlorobenzene-d4	7.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetone	58	C3H6O	000067-64-1	10
3			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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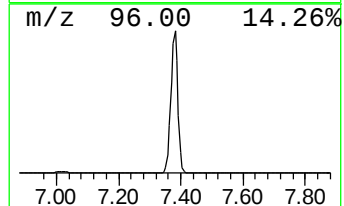
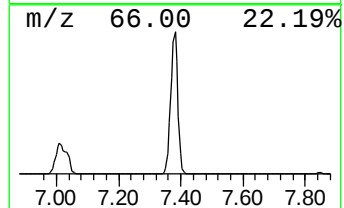
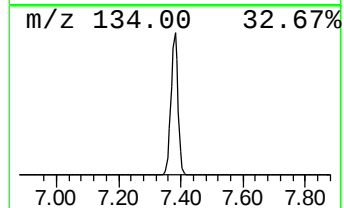
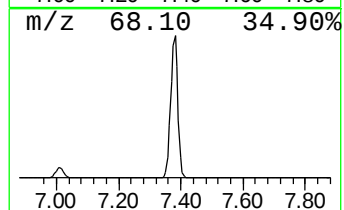
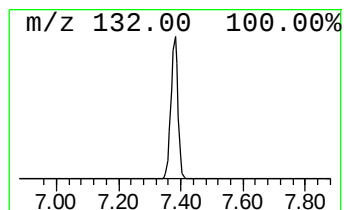
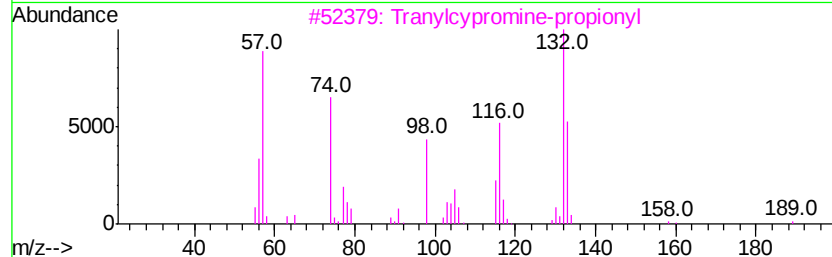
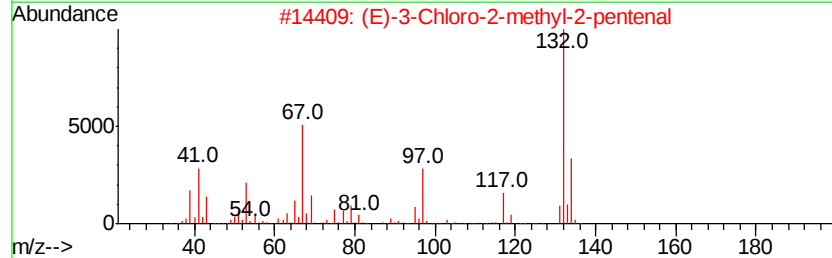
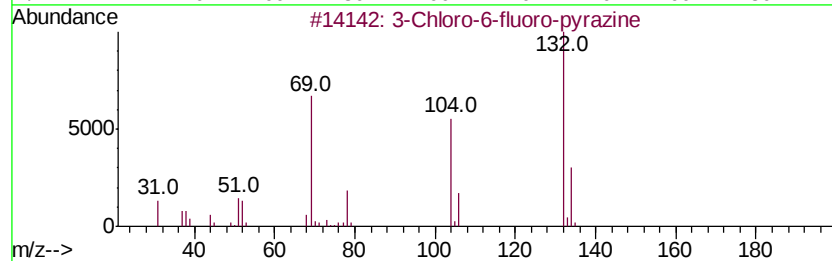
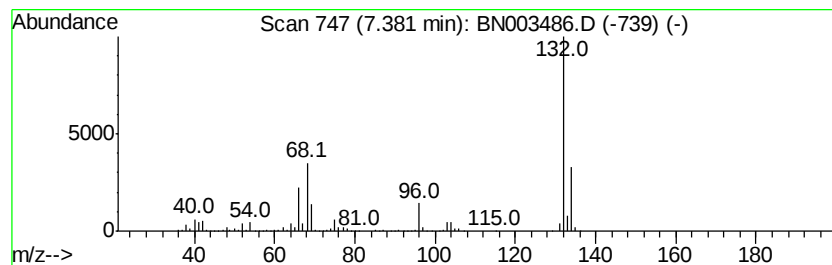
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown7.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	120.27 ng	3390300	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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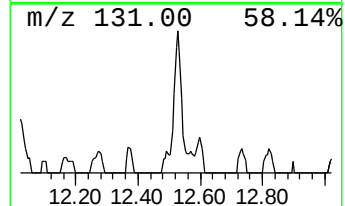
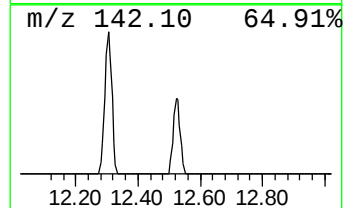
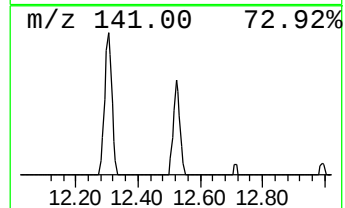
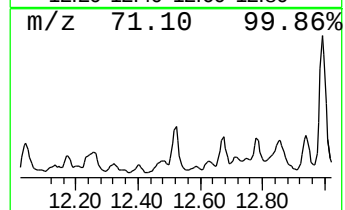
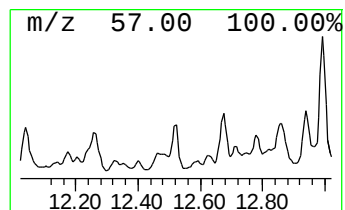
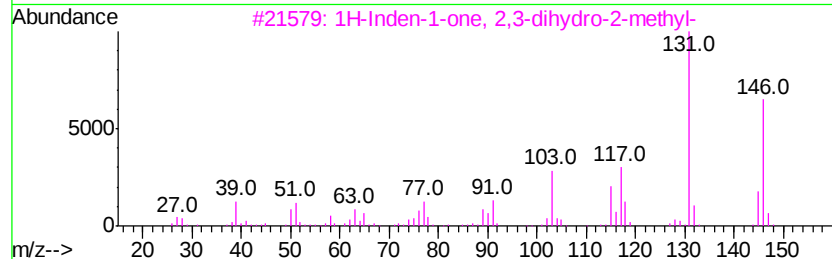
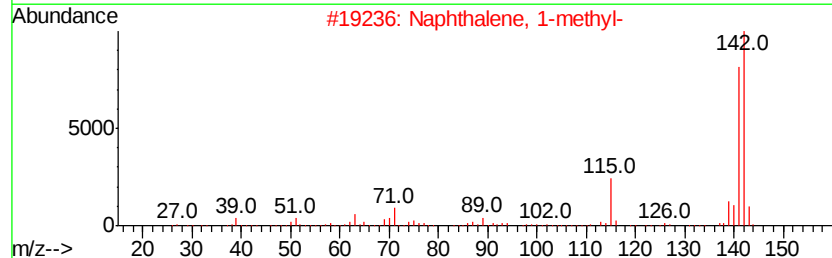
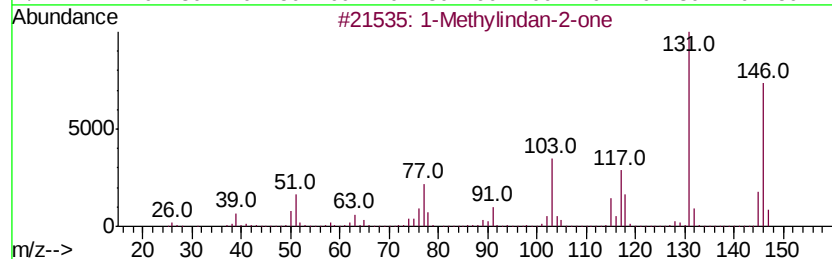
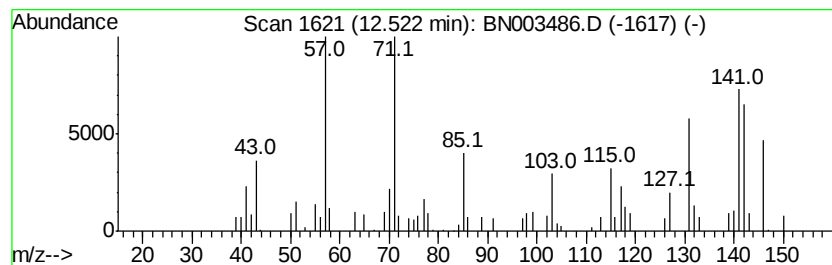
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Peak Number 6 unknown12.52 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	2.13 ng	89098	Naphthalene-d8	10.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylindan-2-one	146	C10H10O	035587-60-1	38
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	35
3		1H-Inden-1-one, 2,3-dihydro-2-methyl-	146	C10H10O	017496-14-9	20
4		Naphthalene, 1,2,3,4-tetrahydro-	146	C11H14	002809-64-5	18
5		Naphthalene, 1,2,3,4-tetrahydro-	146	C11H14	001680-51-9	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
Operator : MJ/SJ
Sample : J5860-21
Misc :
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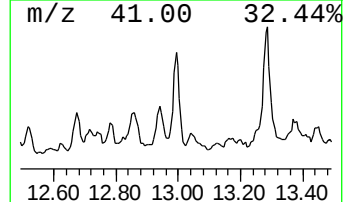
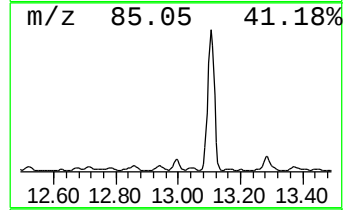
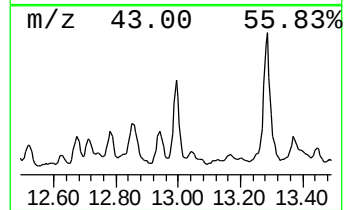
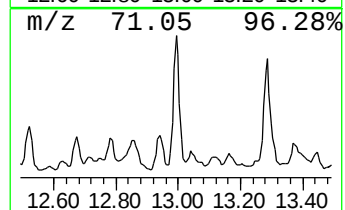
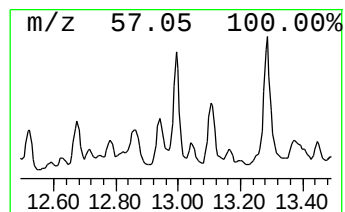
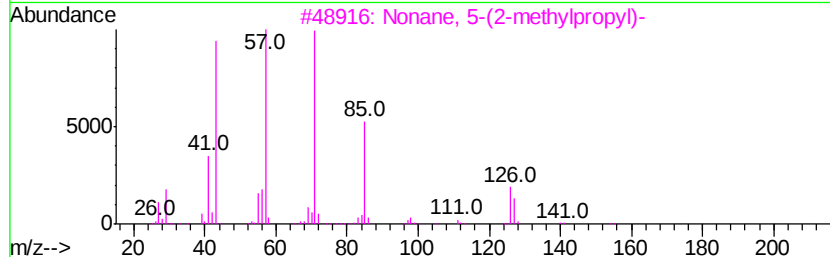
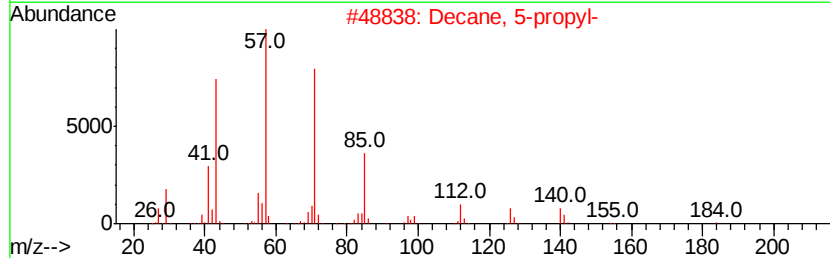
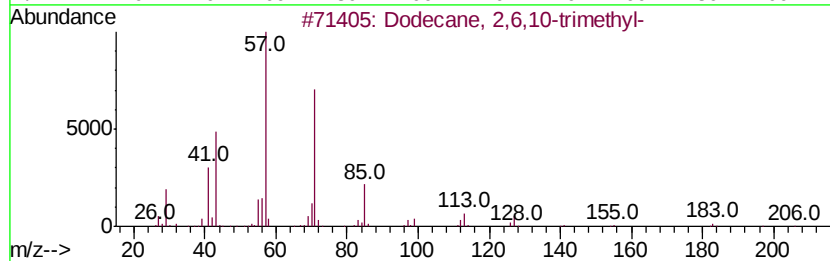
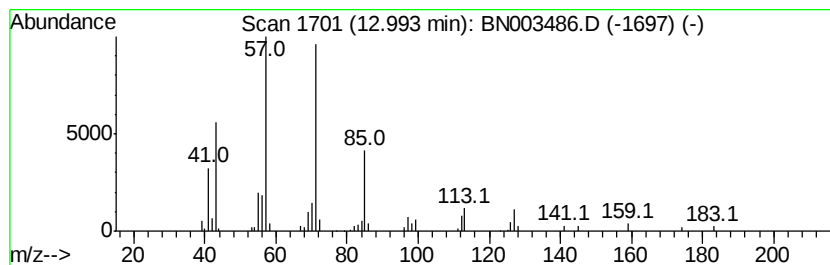
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ClientSampleId :
EP2-SB-110(14-16)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 2,6,10-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	2.01 ng	120702	Acenaphthene-d10	14.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane, 2,6,10-trimethyl-	212 C15H32	003891-98-3	83
2	Decane, 5-propyl-	184 C13H28	017312-62-8	80
3	Nonane, 5-(2-methylpropyl)-	184 C13H28	062185-53-9	64
4	Dodecane, 2,6,11-trimethyl-	212 C15H32	031295-56-4	64
5	Nonane, 5-butyl-	184 C13H28	017312-63-9	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
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Operator : MJ/SJ
Sample : J5860-21
Misc :
ALS Vial : 18 Sample Multiplier: 1

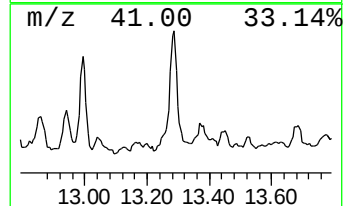
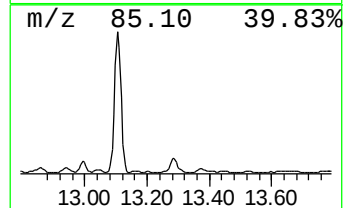
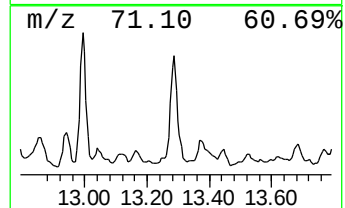
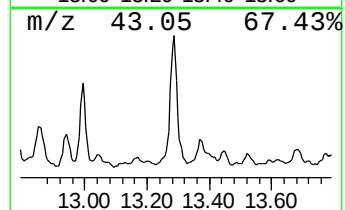
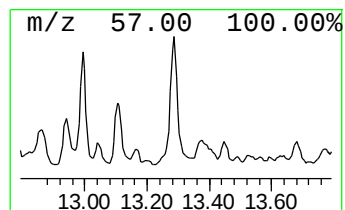
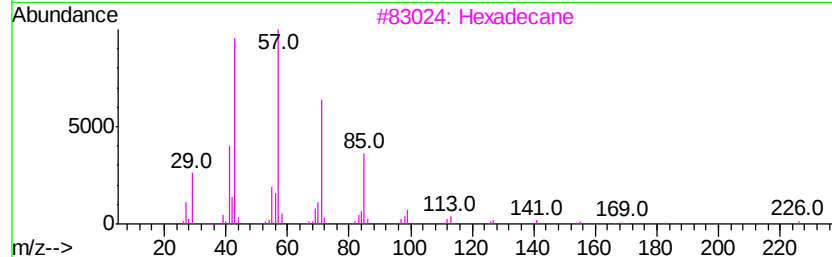
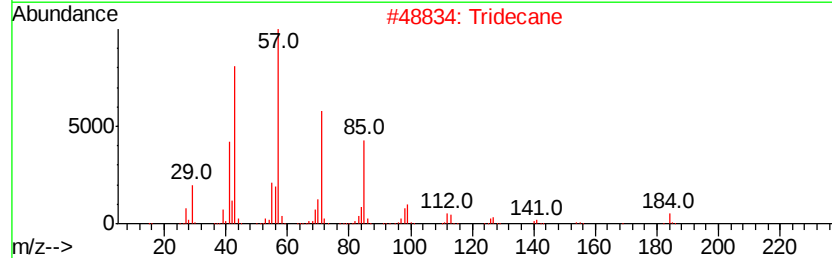
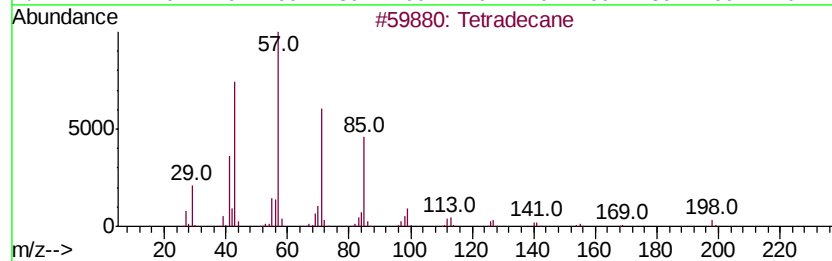
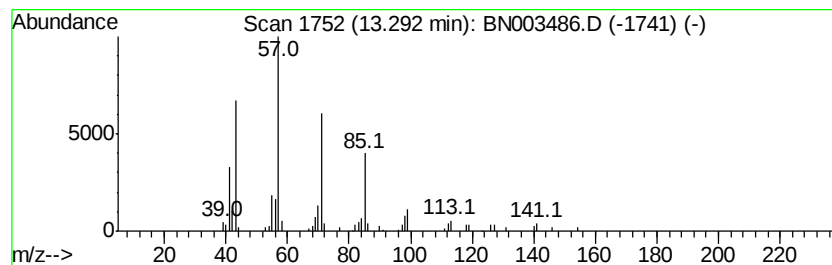
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EP2-SB-110(14-16)

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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 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	3.69 ng	221105	Acenaphthene-d10	14.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecane	198 C14H30	000629-59-4 96
2		Tridecane	184 C13H28	000629-50-5 91
3		Hexadecane	226 C16H34	000544-76-3 91
4		Dodecane	170 C12H26	000112-40-3 90
5		1-Octadecanesulphonyl chloride	352 C18H37ClO2S	1000342-70-4 90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
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Sample : J5860-21
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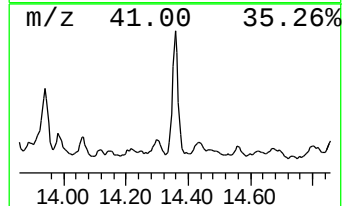
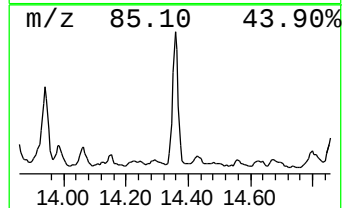
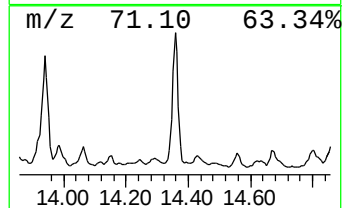
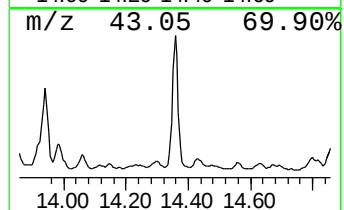
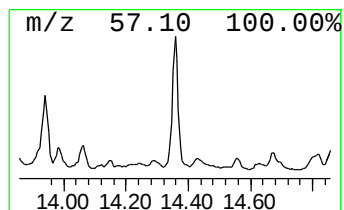
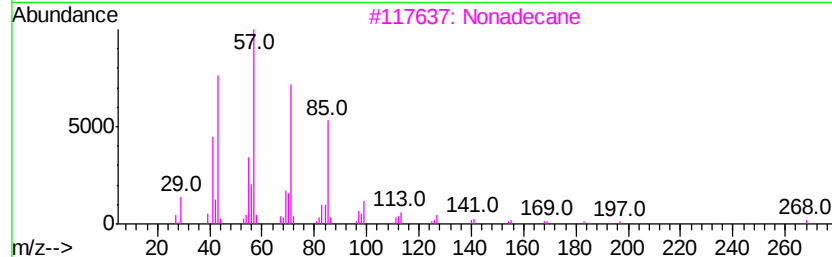
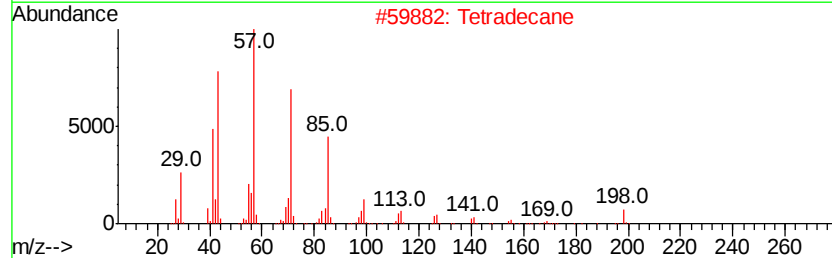
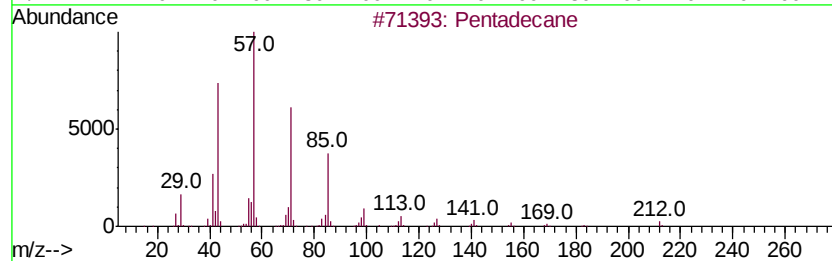
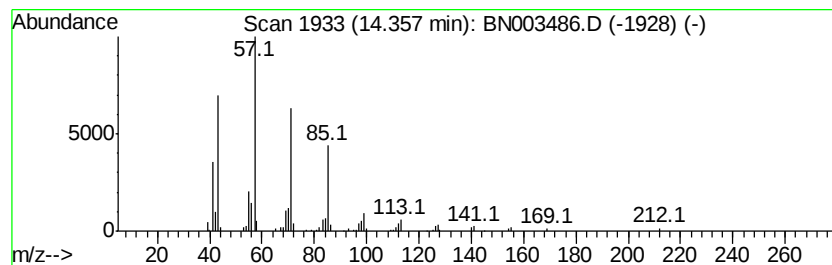
Instrument :
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ClientSampleId :
EP2-SB-110(14-16)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.36	3.84 ng	230231	Acenaphthene-d10		14.49
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	95
2	Tetradecane	198	C14H30	000629-59-4	91
3	Nonadecane	268	C19H40	000629-92-5	86
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5	Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
Operator : MJ/SJ
Sample : J5860-21
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EP2-SB-110(14-16)

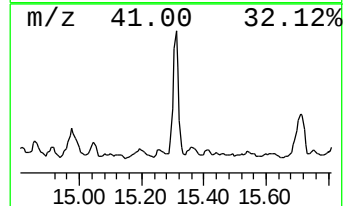
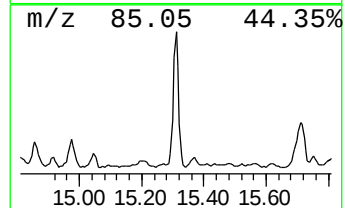
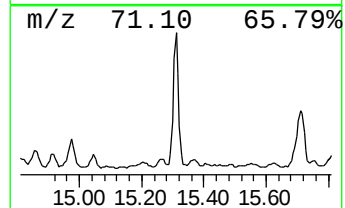
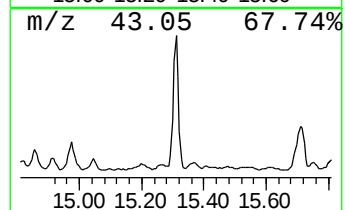
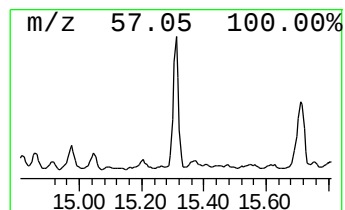
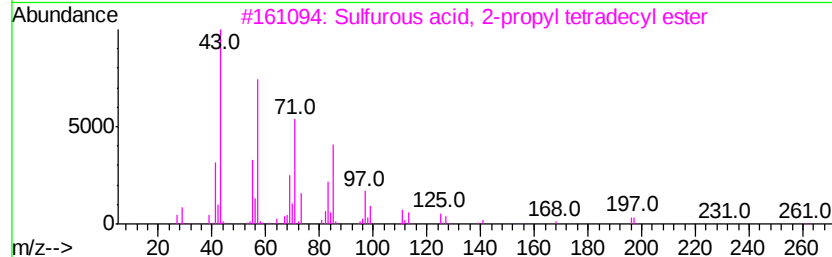
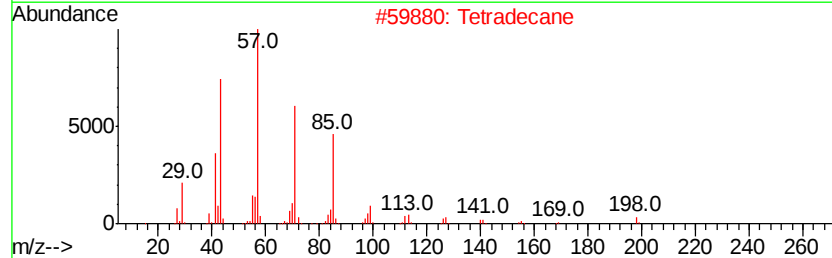
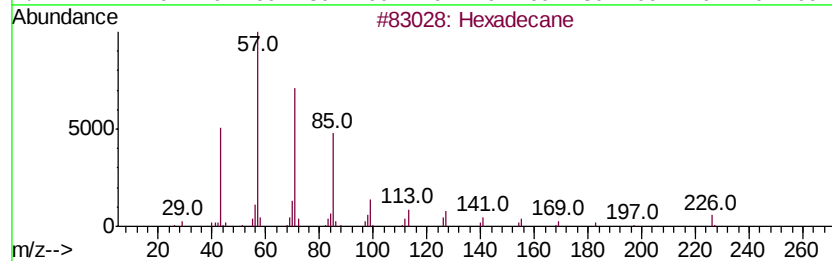
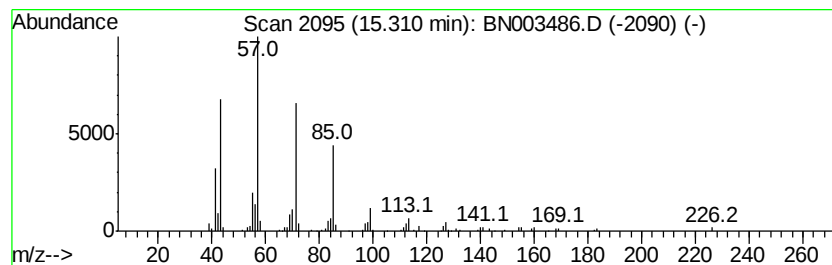
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.31	3.83 ng	229773	Acenaphthene-d10	14.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	97
2			Tetradecane	198	C14H30	000629-59-4	91
3			Sulfurous acid, 2-propyl tetradecyl ester	320	C17H36O3S	1000309-12-5	90
4			Eicosane	282	C20H42	000112-95-8	86
5			Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
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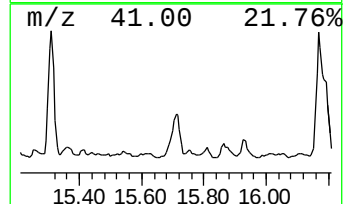
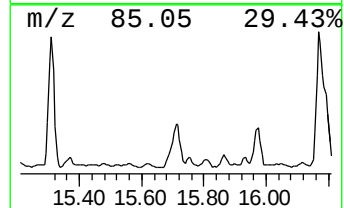
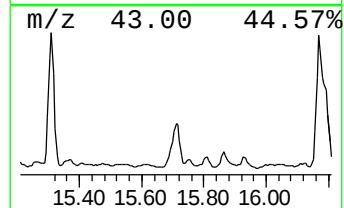
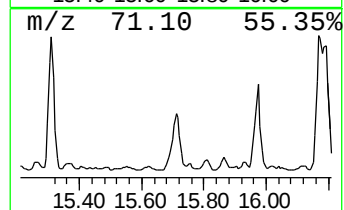
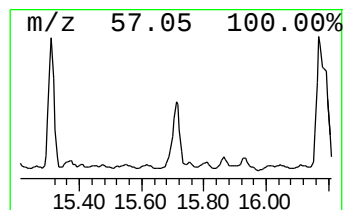
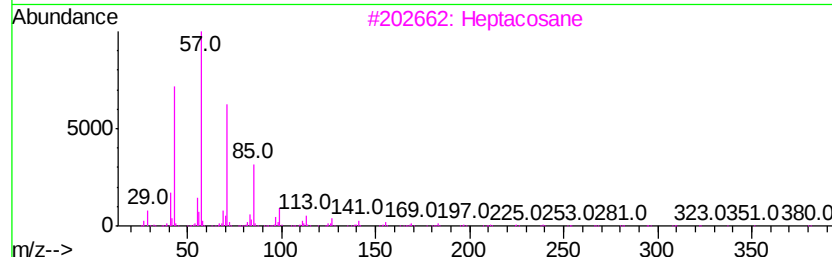
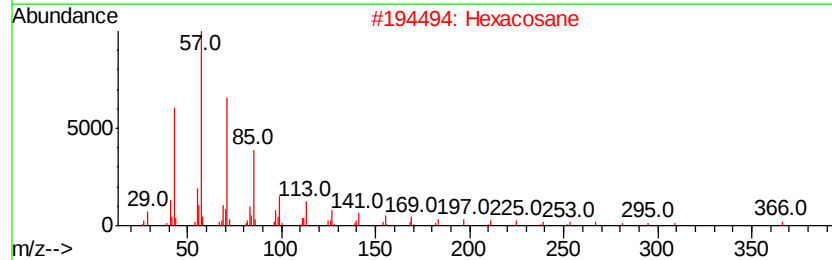
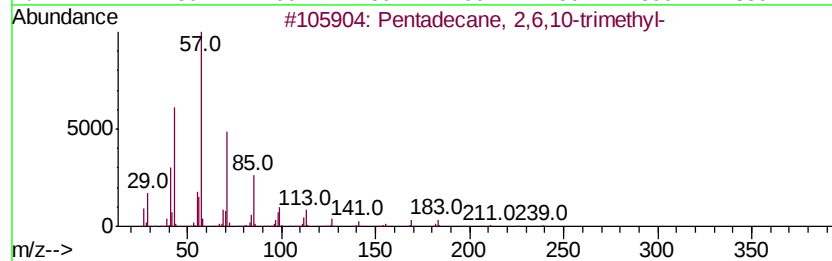
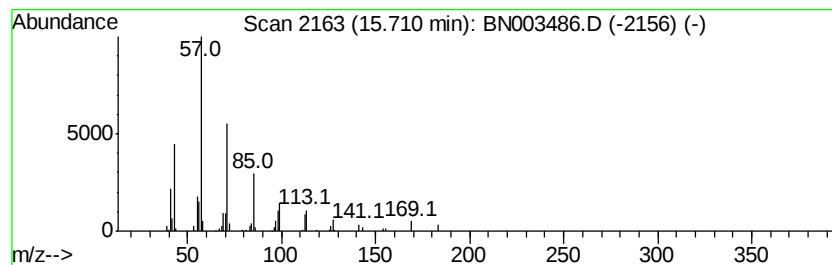
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentadecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	2.48 ng	148384	Acenaphthene-d10	14.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Hexacosane	366	C26H54	000630-01-3	80
3		Heptacosane	380	C27H56	000593-49-7	78
4		Tridecane, 5-propyl-	226	C16H34	055045-11-9	72
5		Tetracosane	338	C24H50	000646-31-1	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
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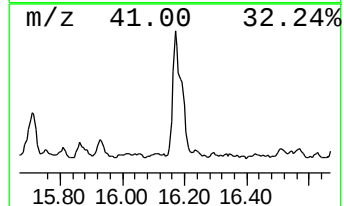
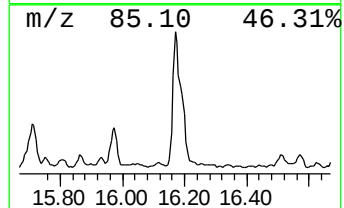
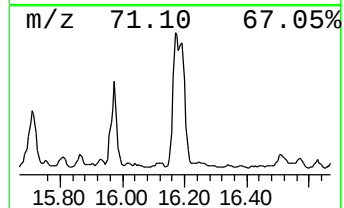
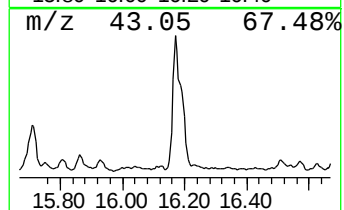
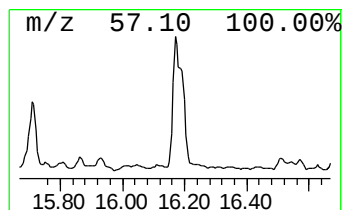
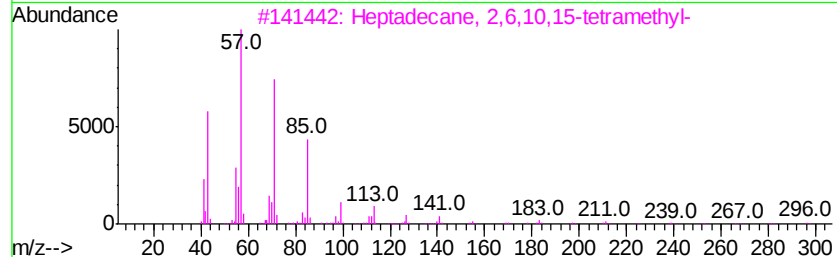
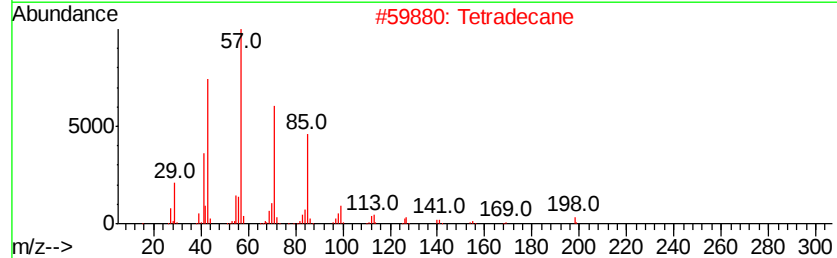
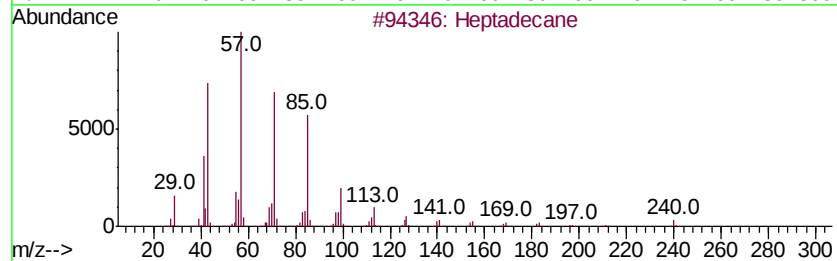
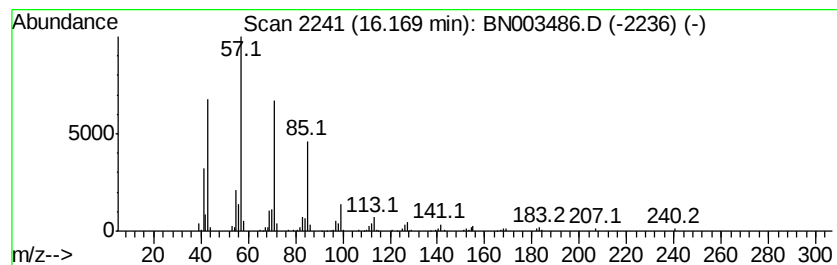
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.17	3.77 ng	262751	Phenanthrene-d10	17.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Tetradecane		198	C14H30	000629-59-4	91
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	91
4	Hexadecane		226	C16H34	000544-76-3	91
5	Pentadecane		212	C15H32	000629-62-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
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ALS Vial : 18 Sample Multiplier: 1

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EP2-SB-110(14-16)

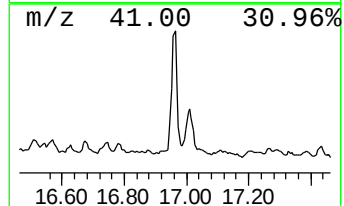
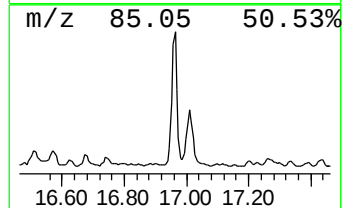
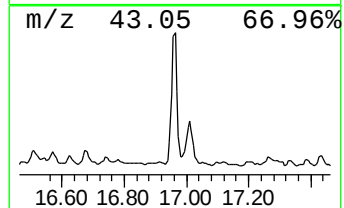
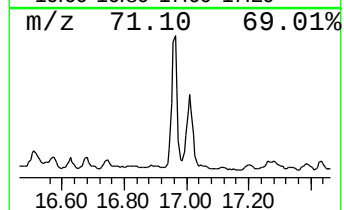
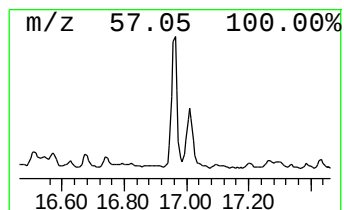
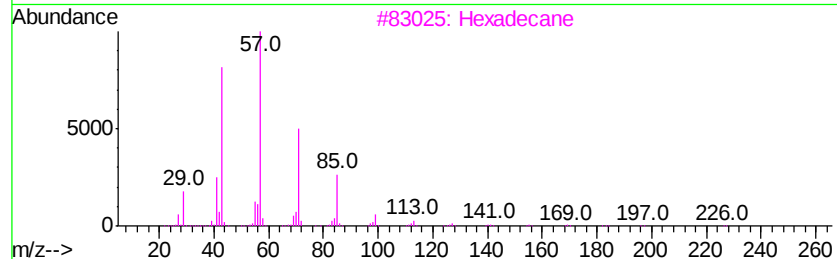
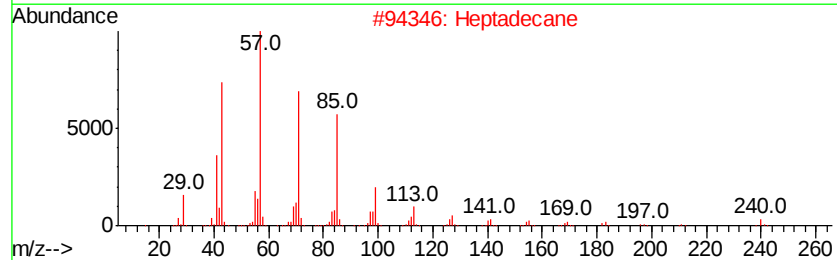
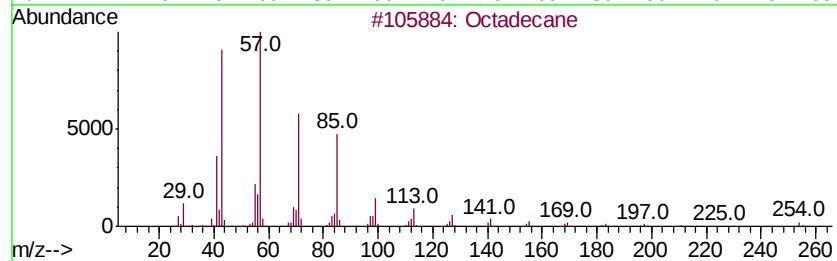
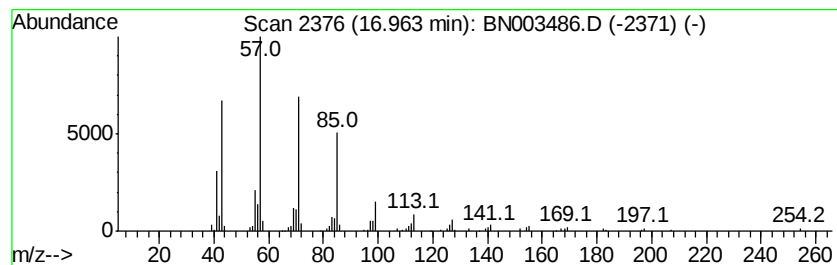
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.96	2.58 ng	179365	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	92
2		Heptadecane	240	C17H36	000629-78-7	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Heneicosane	296	C21H44	000629-94-7	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
Operator : MJ/SJ
Sample : J5860-21
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EP2-SB-110(14-16)

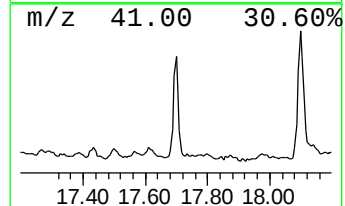
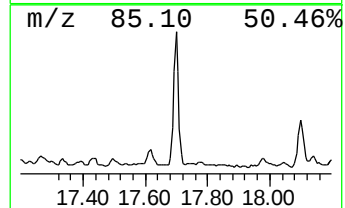
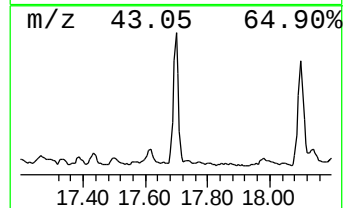
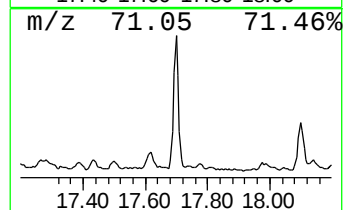
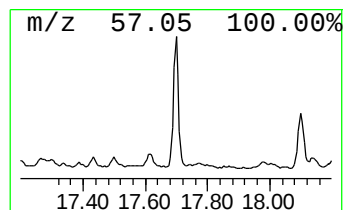
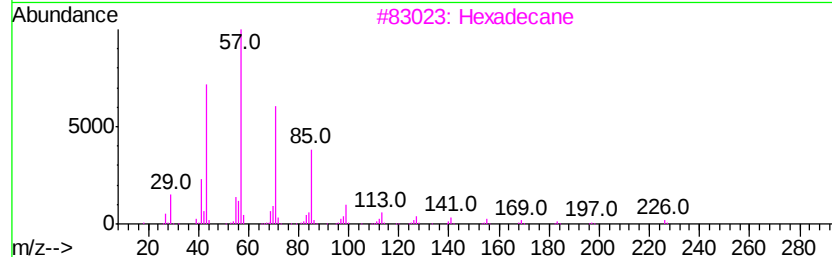
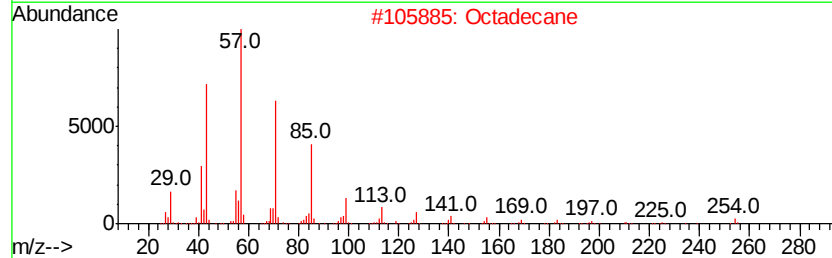
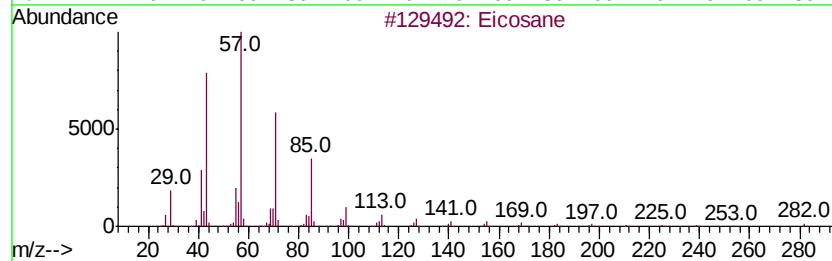
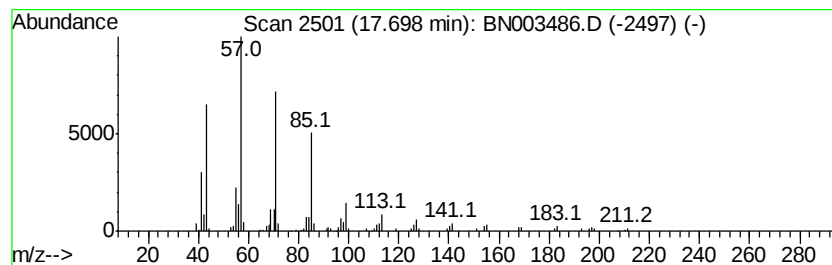
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.70	2.19 ng	152230	Phenanthrene-d10	17.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Octadecane	254	C18H38	000593-45-3	91
3			Hexadecane	226	C16H34	000544-76-3	91
4			Heptadecane	240	C17H36	000629-78-7	91
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
Operator : MJ/SJ
Sample : J5860-21
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EP2-SB-110(14-16)

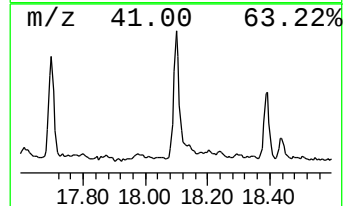
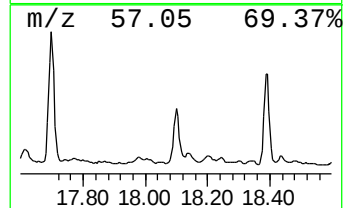
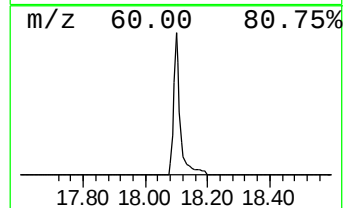
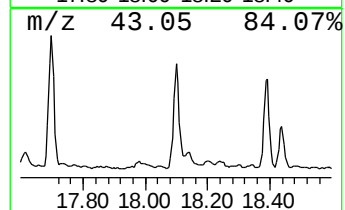
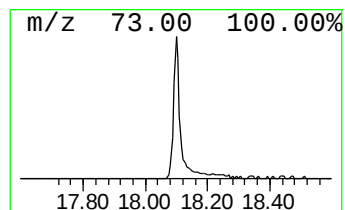
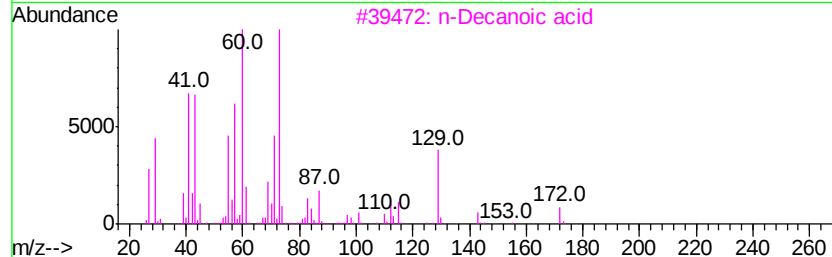
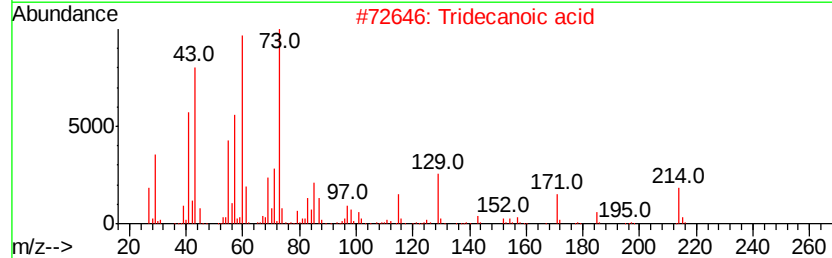
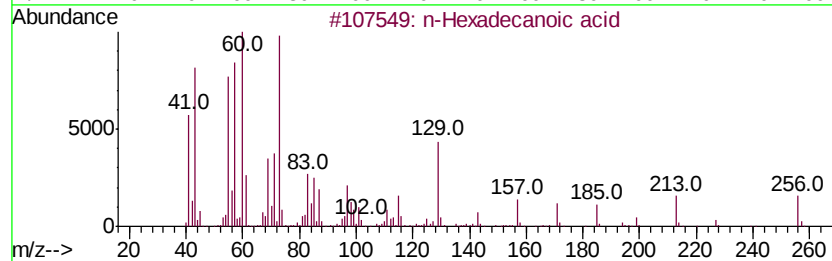
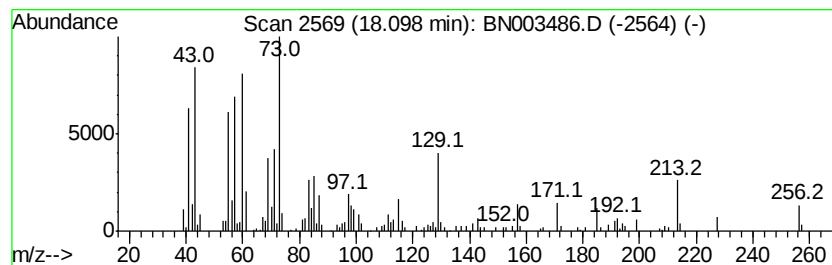
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.10	3.69 ng	256642	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		n-Decanoic acid	172	C10H20O2	000334-48-5	76
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110918\
Data File : BN003486.D
Acq On : 09 Nov 2018 23:27
Operator : MJ/SJ
Sample : J5860-21
Misc :
ALS Vial : 18 Sample Multiplier: 1

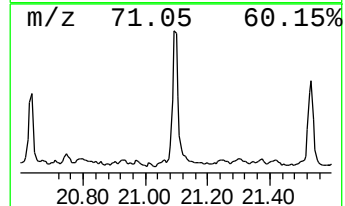
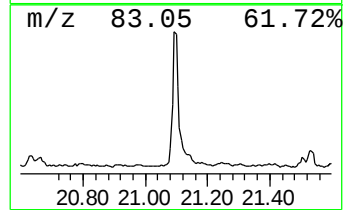
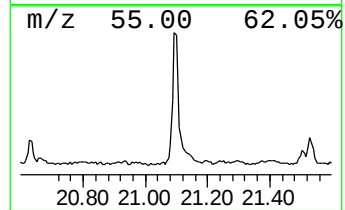
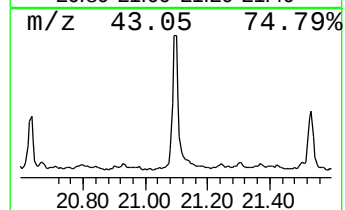
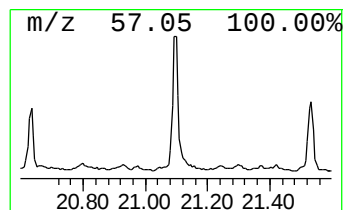
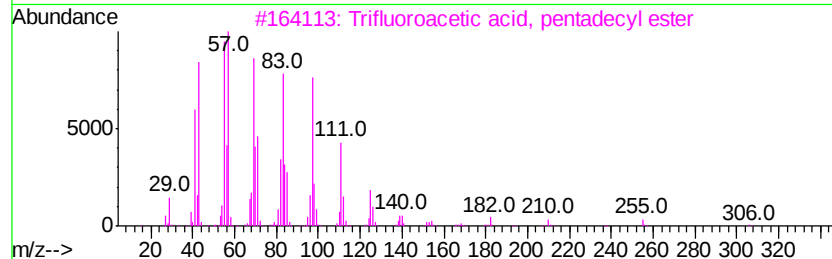
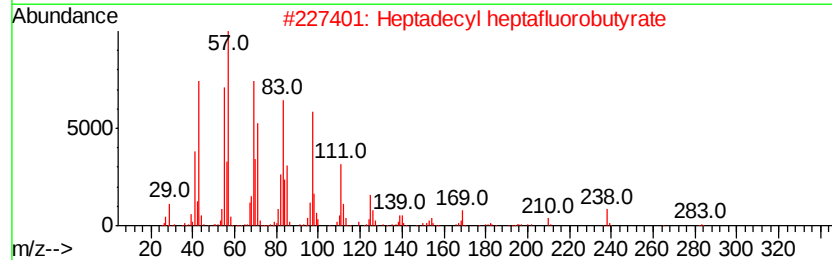
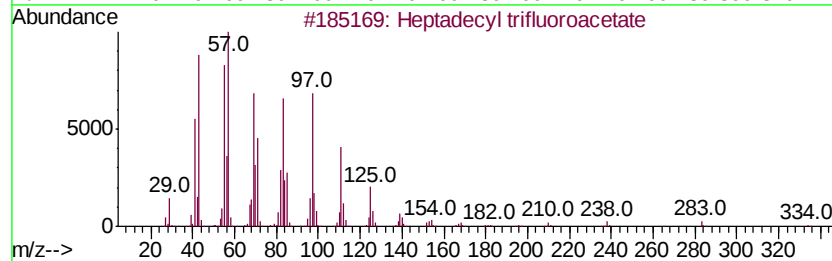
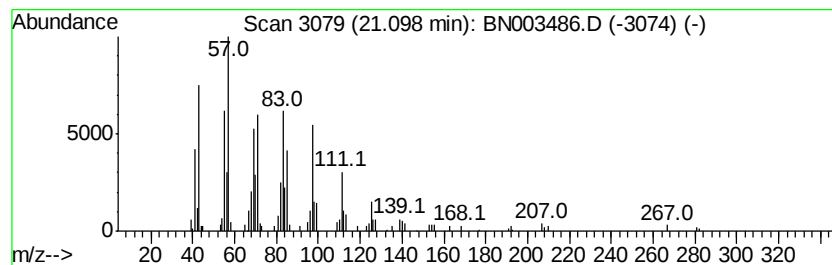
Instrument :
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ClientSampleId :
EP2-SB-110(14-16)

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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 Heptadecyl trifluoroacetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.10	2.63 ng	166553	Chrysene-d12	21.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
3	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110918\
 Data File : BN003486.D
 Acq On : 09 Nov 2018 23:27
 Operator : MJ/SJ
 Sample : J5860-21
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EP2-SB-110(14-16)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.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
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unknown3.56	3.56	2.4	ng	68741	1	7.85	563779	20.0
2-Pentanone, 4-hy...	4.96	3.9	ng	108983	1	7.85	563779	20.0
unknown7.38	7.38	120.3	ng	3390300	1	7.85	563779	20.0
unknown12.52	12.52	2.1	ng	89098	2	10.65	835545	20.0
Dodecane, 2,6,10-...	12.99	2.0	ng	120702	3	14.49	1198740	20.0
Tetradecane	13.29	3.7	ng	221105	3	14.49	1198740	20.0
Pentadecane	14.36	3.8	ng	230231	3	14.49	1198740	20.0
Hexadecane	15.31	3.8	ng	229773	3	14.49	1198740	20.0
Pentadecane, 2,6,...	15.71	2.5	ng	148384	3	14.49	1198740	20.0
Heptadecane	16.17	3.8	ng	262751	4	17.23	1392710	20.0
Octadecane	16.96	2.6	ng	179365	4	17.23	1392710	20.0
Eicosane	17.70	2.2	ng	152230	4	17.23	1392710	20.0
n-Hexadecanoic acid	18.10	3.7	ng	256642	4	17.23	1392710	20.0
Heptadecyl triflu...	21.10	2.6	ng	166553	5	21.40	1268950	20.0