

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
 Data File : BF077935.D
 Acq On : 25 Mar 2015 4:35
 Operator : TP/IZ
 Sample : G1613-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4-1

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-BF031115.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	1.047	11	14	18	rVB	69679	122185	11.00%	1.150%
2	1.275	31	34	36	rVB	23822	34311	3.09%	0.323%
3	1.390	41	44	47	rVB	105148	162458	14.62%	1.529%
4	1.561	56	59	65	rVB	38067	56126	5.05%	0.528%
5	1.653	65	67	71	rVV	13120	17661	1.59%	0.166%
6	1.767	75	77	89	rVB	85117	143181	12.89%	1.348%
7	4.864	346	348	355	rVB	87209	109130	9.82%	1.027%
8	5.402	393	395	398	rBV	594817	695848	62.63%	6.551%
9	6.693	506	508	511	rBV	674233	702447	63.23%	6.613%
10	6.830	518	520	523	rBV	856529	779421	70.16%	7.337%
11	7.116	543	545	548	rBV	283066	250480	22.55%	2.358%
12	7.299	559	561	564	rBV	573596	598772	53.90%	5.637%
13	7.813	603	606	608	rBV	518858	462930	41.67%	4.358%
14	8.693	681	683	686	rBV	407807	350757	31.57%	3.302%
15	10.042	799	801	803	rBV	1101500	959231	86.34%	9.030%
16	10.545	843	845	848	rVB	19875	25188	2.27%	0.237%
17	10.865	871	873	875	rBV	426925	406097	36.55%	3.823%
18	11.837	956	958	961	rBV2	388499	520575	46.86%	4.901%
19	12.042	974	976	979	rBV	14200	16682	1.50%	0.157%
20	12.602	1023	1025	1027	rBV	13531	14070	1.27%	0.132%
21	12.694	1031	1033	1036	rBV	294987	414471	37.31%	3.902%
22	13.128	1070	1071	1073	rBV	10272	12882	1.16%	0.121%
23	14.694	1205	1208	1210	rBV	1101313	1110971	100.00%	10.459%
24	15.037	1235	1238	1240	rBV2	7840	12558	1.13%	0.118%
25	15.334	1261	1264	1266	rBV	11719	15480	1.39%	0.146%
26	15.414	1268	1271	1273	rBV3	10179	20741	1.87%	0.195%
27	15.883	1310	1312	1318	rBV	43916	94323	8.49%	0.888%
28	15.974	1318	1320	1323	rBV	369564	489929	44.10%	4.612%
29	16.146	1333	1335	1337	rBV	24313	30196	2.72%	0.284%
30	16.283	1345	1347	1350	rBV4	12047	12929	1.16%	0.122%
31	16.637	1375	1378	1380	rBV2	22941	34805	3.13%	0.328%
32	16.683	1380	1382	1384	rBV	17046	23010	2.07%	0.217%
33	16.946	1404	1405	1409	rVB4	13169	18027	1.62%	0.170%
34	17.277	1429	1434	1435	rBV4	24450	69865	6.29%	0.658%

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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

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.471	1449	1451	1453	rBV2	19363	31026	2.79%	0.292%
36	17.723	1470	1473	1476	rBV	403712	510482	45.95%	4.806%
37	19.266	1604	1608	1615	rVB	433378	862700	77.65%	8.121%
38	19.574	1631	1635	1639	rVB2	80531	185936	16.74%	1.750%
39	20.020	1670	1674	1680	rBV4	65927	163619	14.73%	1.540%
40	20.409	1705	1708	1716	rVB10	28113	81058	7.30%	0.763%

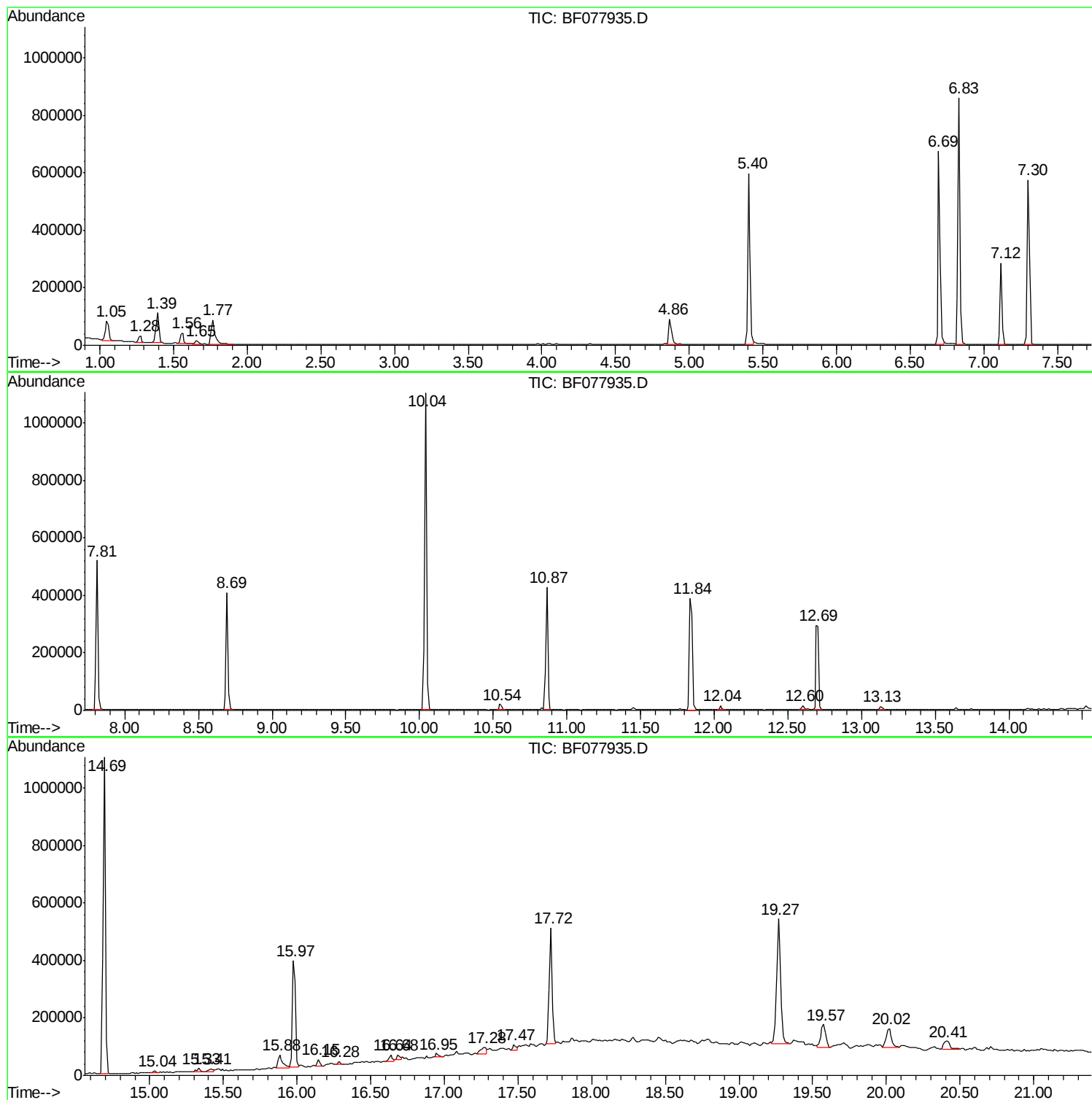
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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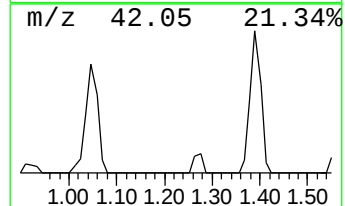
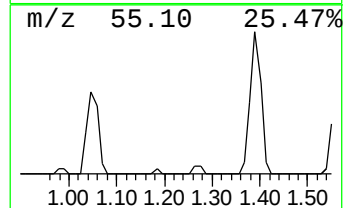
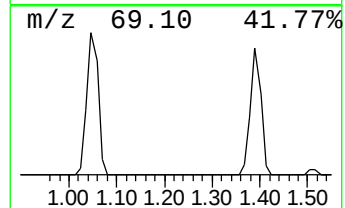
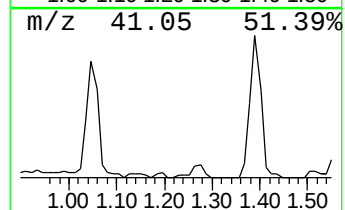
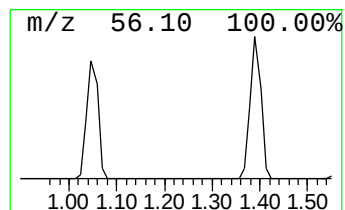
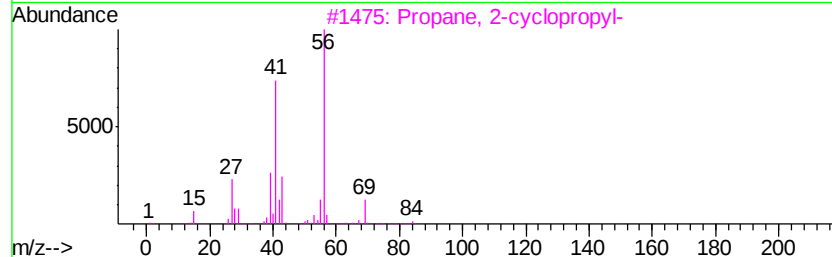
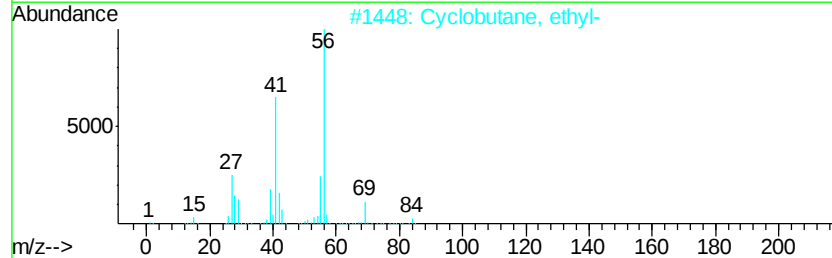
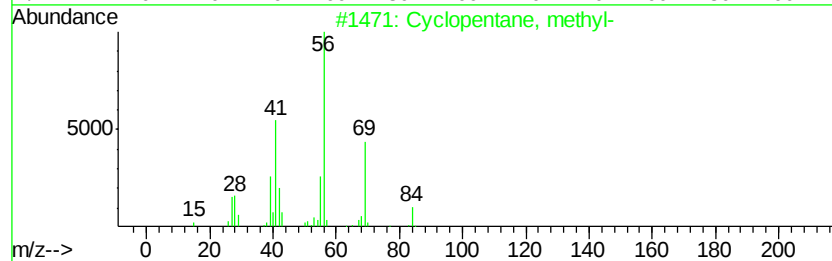
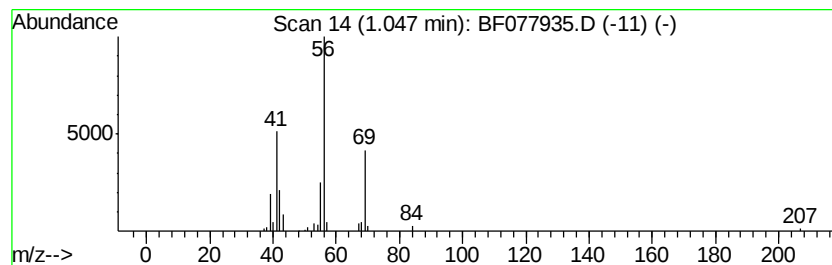
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.05	9.76 ng	122185	1,4-Dichlorobenzene-d4	7.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	72
4		Cyclohexane	84	C6H12	000110-82-7	56
5		Heptane, 4-methylene-	112	C8H16	015918-08-8	56



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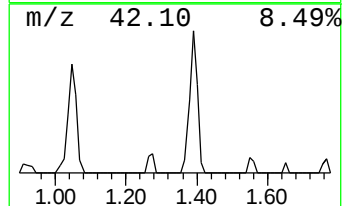
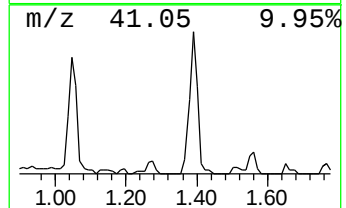
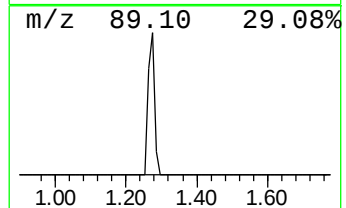
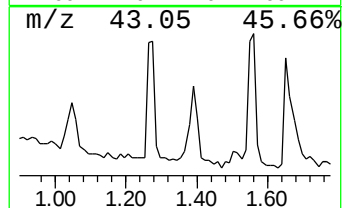
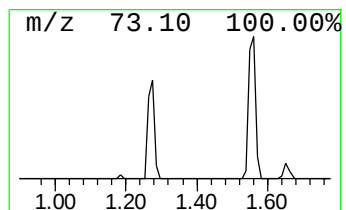
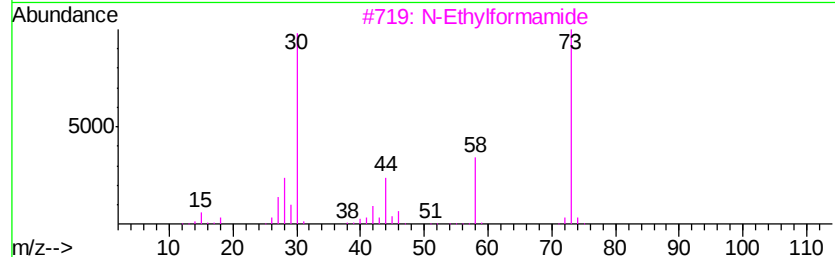
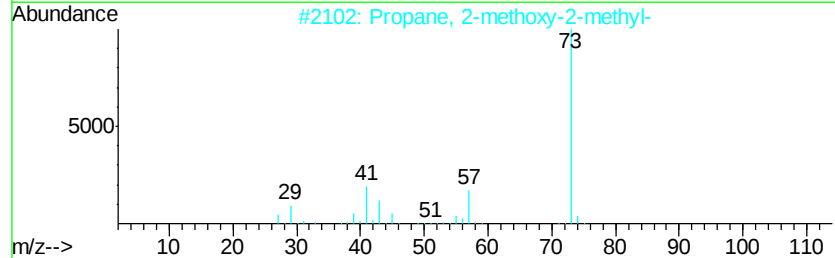
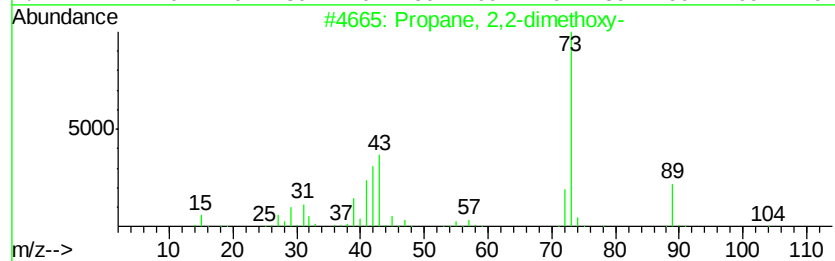
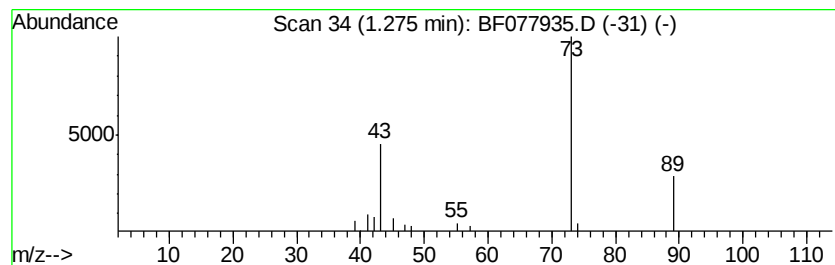
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Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.28	2.74 ng	34311	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	78
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		N-Ethylformamide	73	C3H7NO	000627-45-2	7
4		.beta.-Alanine	89	C3H7NO2	000107-95-9	4
5		Isobutylamine	73	C4H11N	000078-81-9	4



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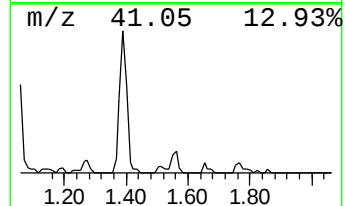
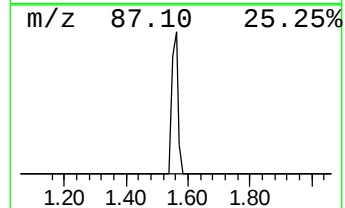
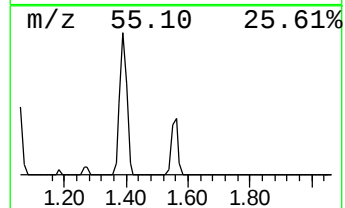
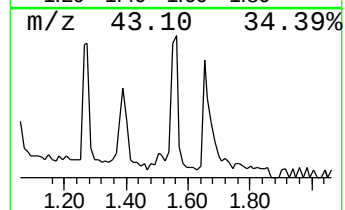
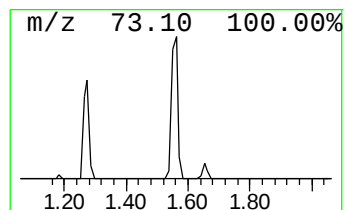
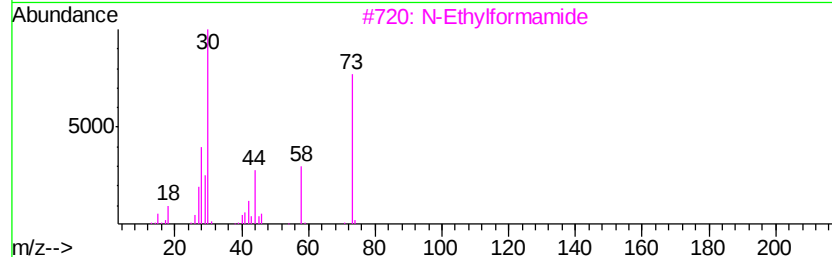
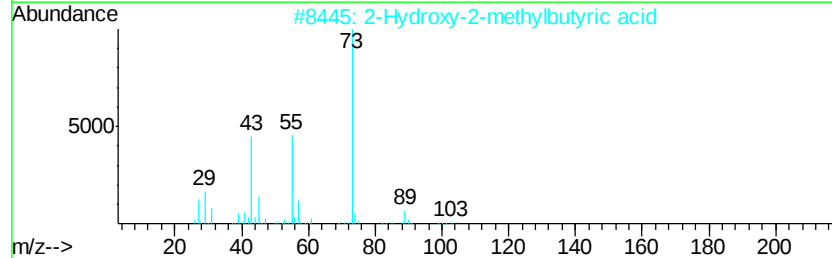
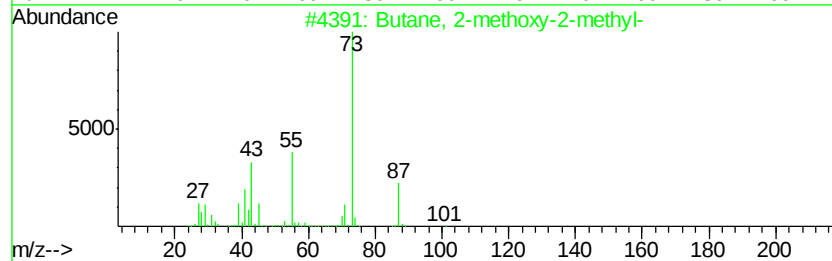
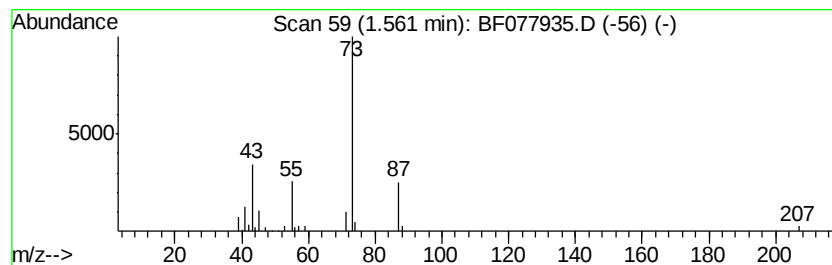
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.56	4.48 ng	56126	1,4-Dichlorobenzene-d4	7.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	12
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9
5		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9



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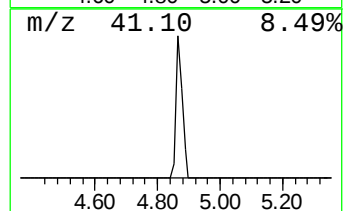
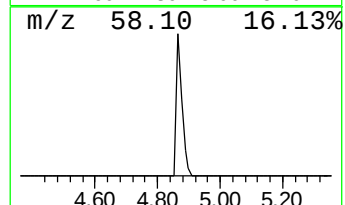
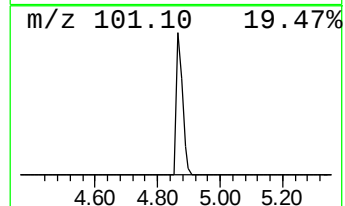
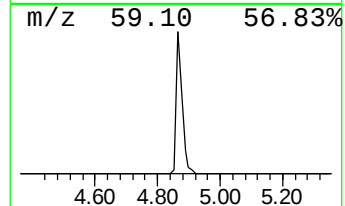
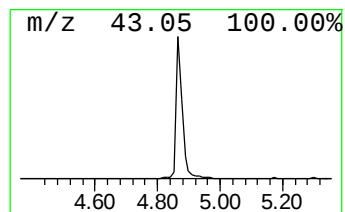
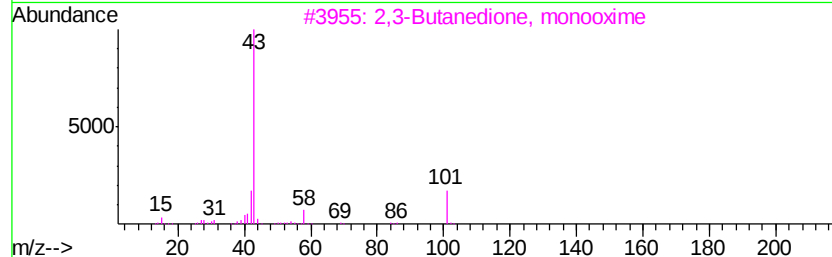
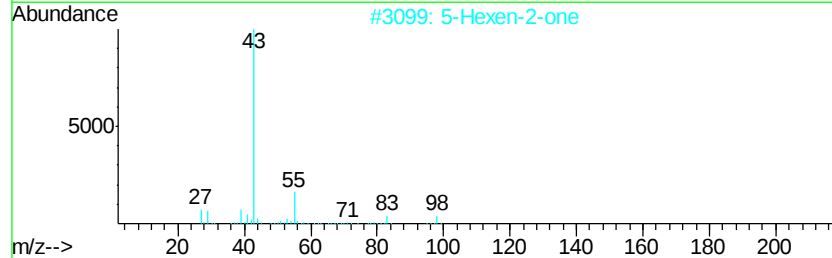
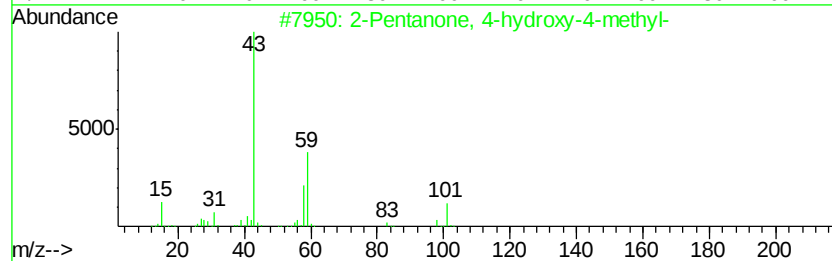
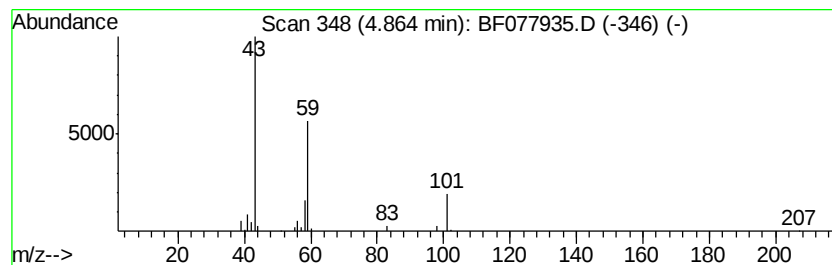
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	8.71 ng	109130	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2,5,8,11,14-Pentaoxapentadecane	222	C10H22O5	000143-24-8	9



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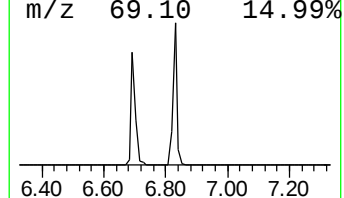
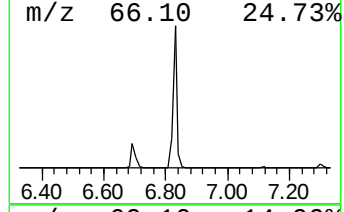
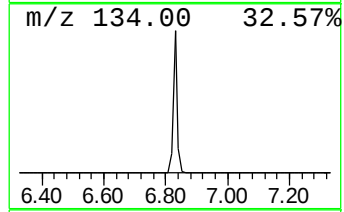
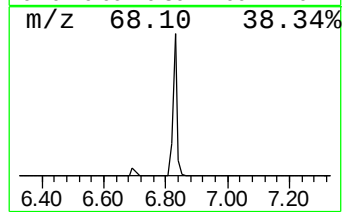
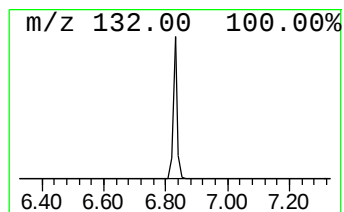
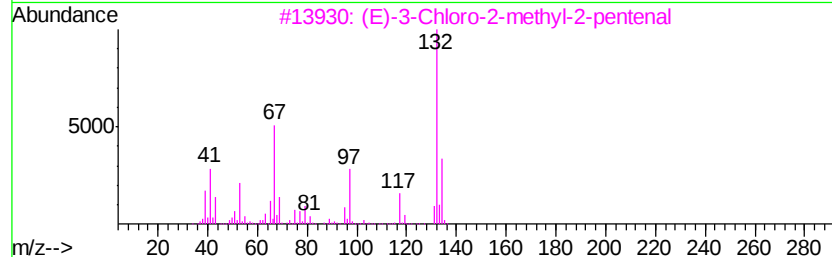
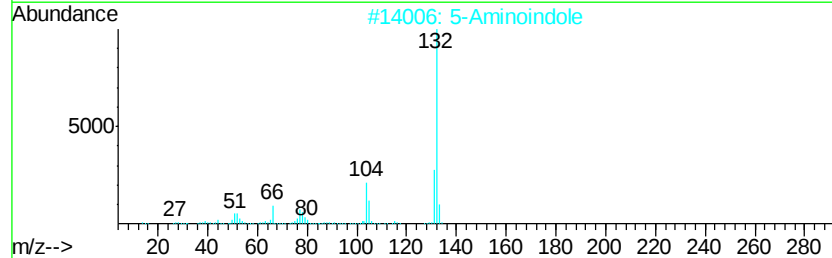
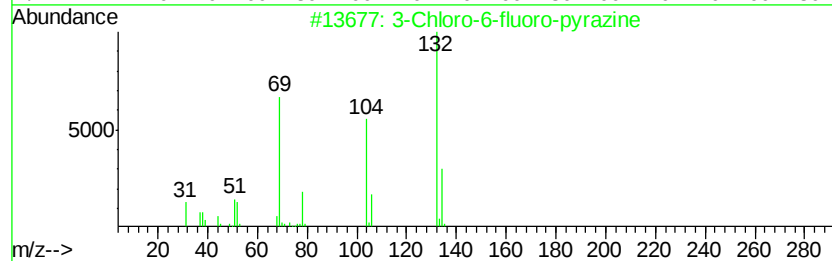
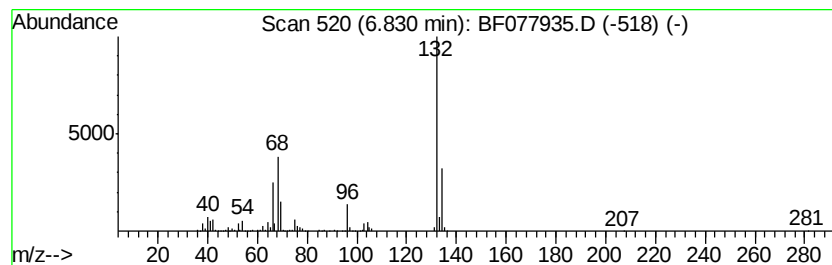
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Peak Number 7 unknown6.83 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	62.23 ng	779421	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

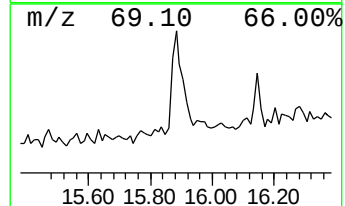
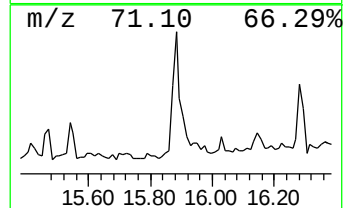
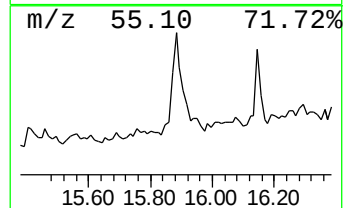
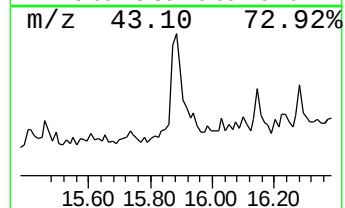
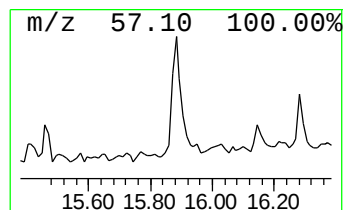
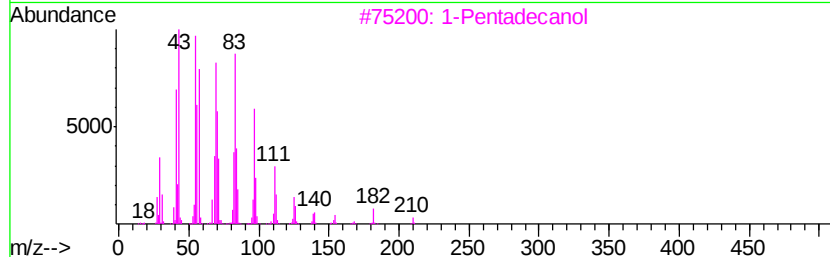
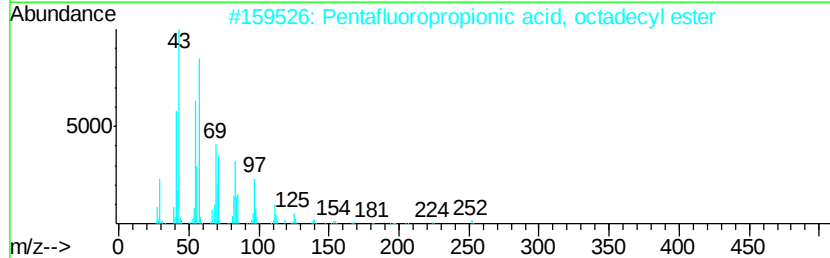
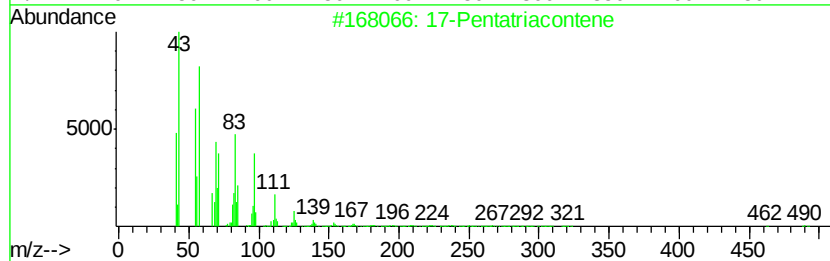
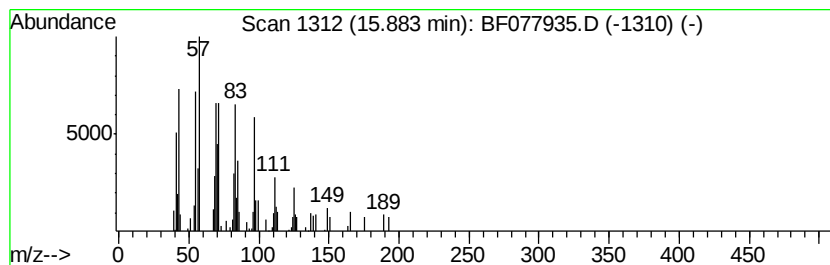
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ClientSampleId :
SB-4-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 17-Pentatriacontene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	3.85 ng	94323	Chrysene-d12	15.97
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	17-Pentatriacontene	491 C35H70	006971-40-0	68
2	Pentafluoropropionic acid, octad...	416 C21H37F5O2	1000280-07-7	64
3	1-Pentadecanol	228 C15H32O	000629-76-5	58
4	Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1	58
5	1-Hentetracontanol	593 C41H84O	040710-42-7	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-4-1

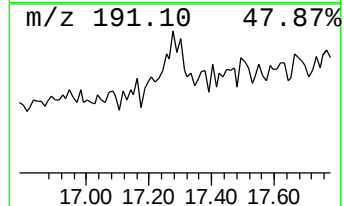
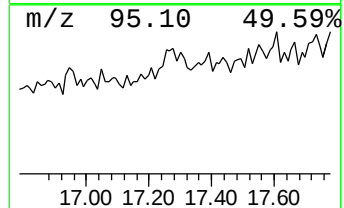
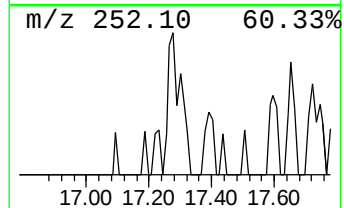
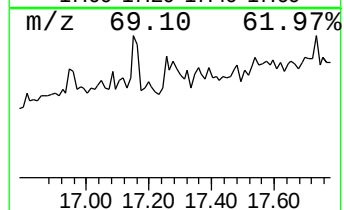
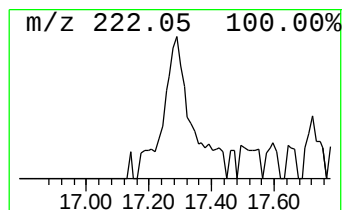
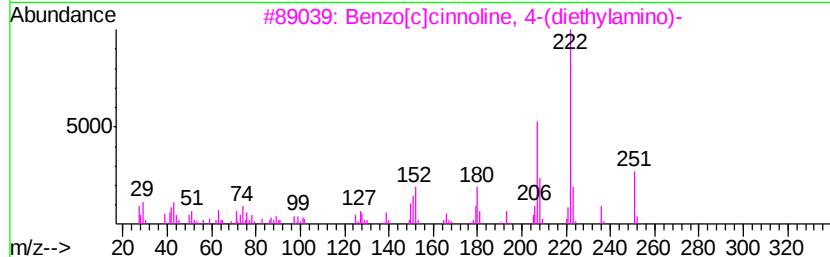
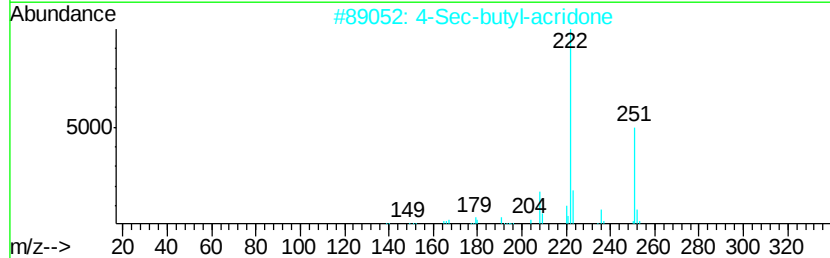
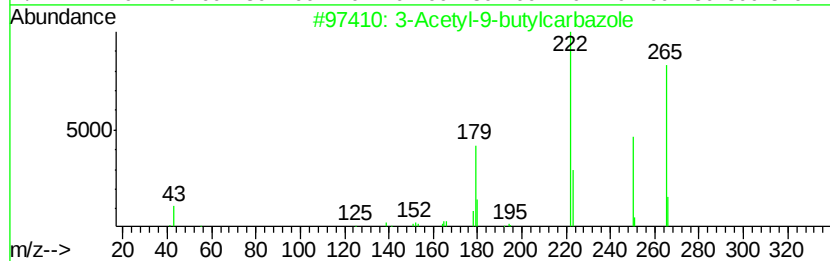
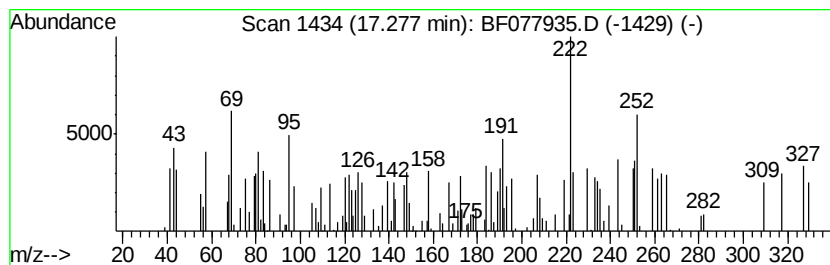
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown17.28 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.28	2.74 ng	69865	Perylene-d12	17.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Acetyl-9-butylcarbazole	265	C18H19NO	001484-01-1	25
2		4-Sec-butyl-acridone	251	C17H17NO	055751-70-7	22
3		Benzo[c]cinnoline, 4-(diethylami...	251	C16H17N3	019174-76-6	22
4		1,3-Dimethyl-2,2-diphenylaziridine	223	C16H17N	1000188-89-9	15
5		trans-2,4-Dimethoxy-5-ethoxy-.be...	222	C13H18O3	1000122-80-1	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

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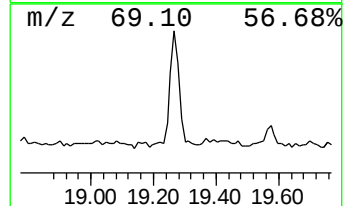
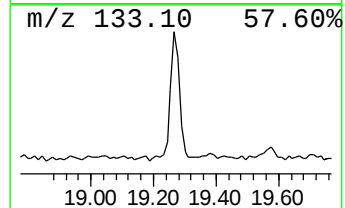
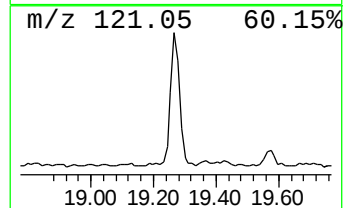
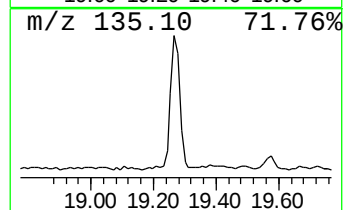
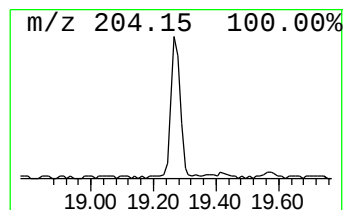
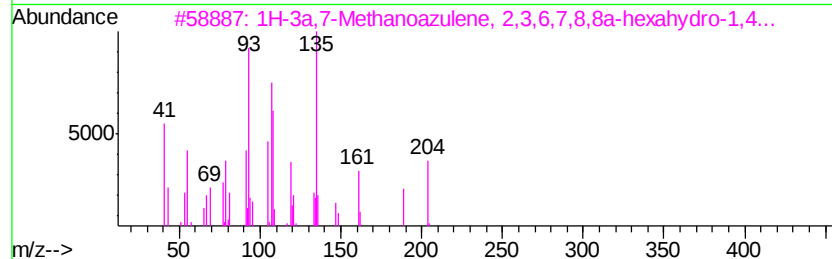
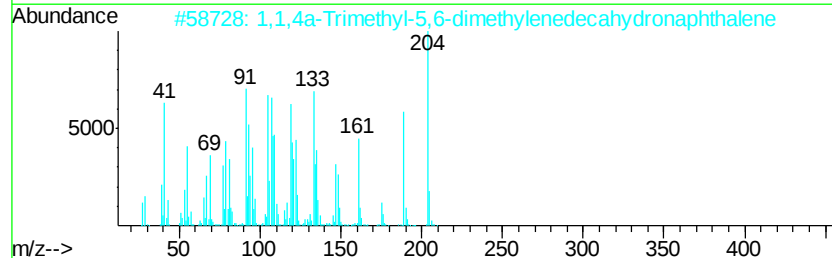
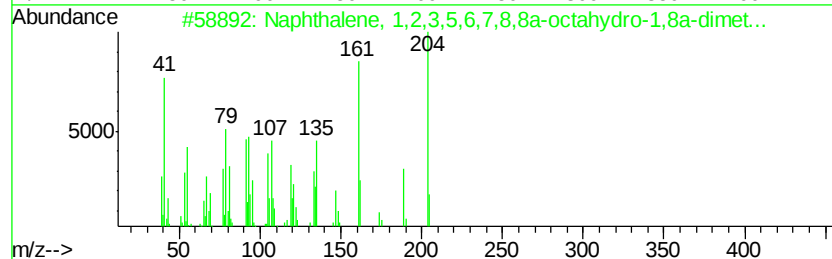
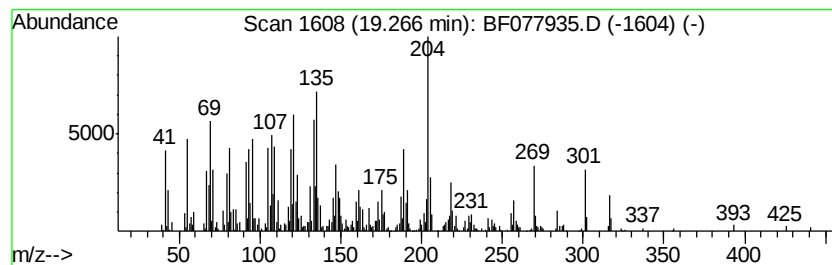
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,2,3,5,6,7,8,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	33.80 ng	862700	Perylene-d12	17.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	59
2		1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	58
3		1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	53
4		Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	53
5		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-4-1

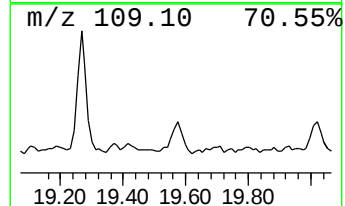
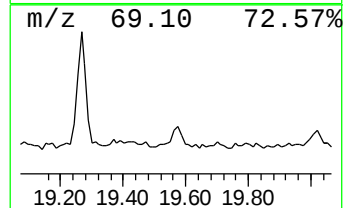
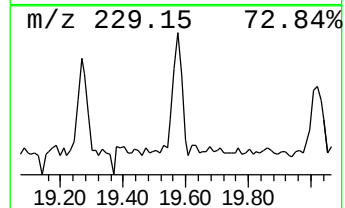
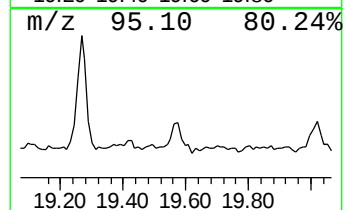
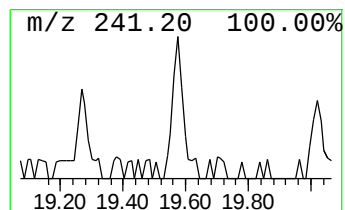
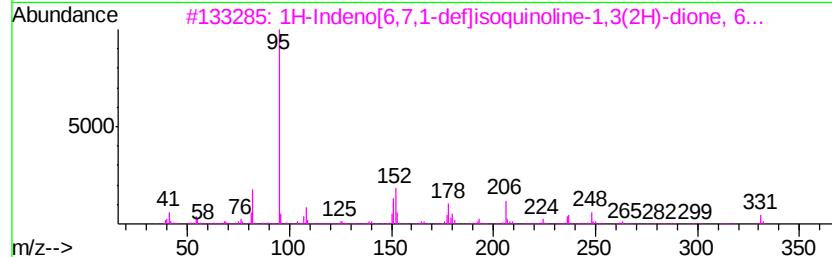
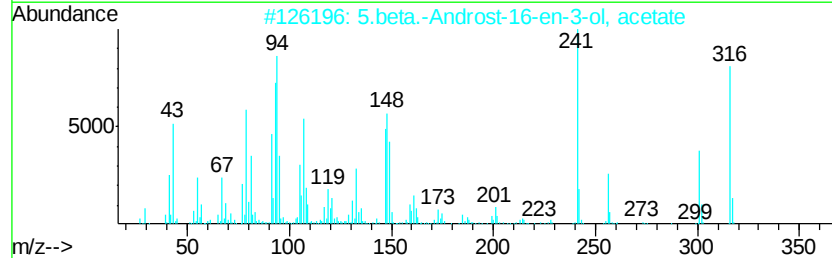
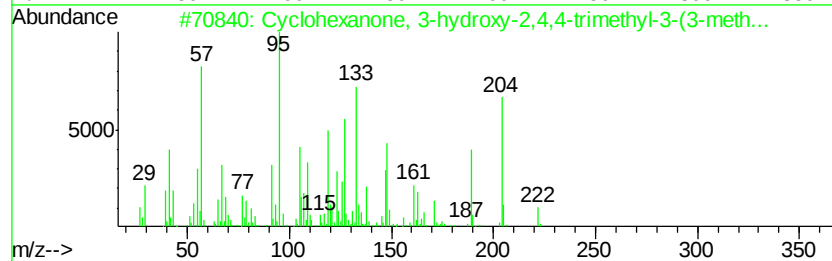
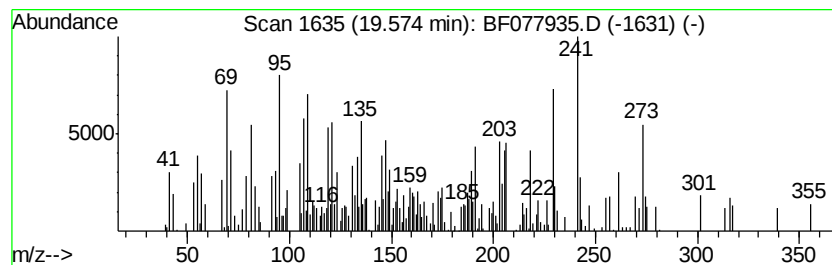
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown19.57 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.57	7.28 ng	185936	Perylene-d12	17.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3-hydroxy-2,4,4-t...	222	C14H22O2	088764-84-5	15
2		5.beta.-Androst-16-en-3-ol, acetate	316	C21H32O2	1000196-50-8	11
3		1H-Indeno[6,7,1-def]isoquinoline...	331	C20H17N3O2	1000272-50-2	10
4		3a,6-Epoxy-3aH-isoindole, 1,2,3,...	257	C16H19NO2	071840-22-7	10
5		Longifolenaldehyde	220	C15H24O	019890-84-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

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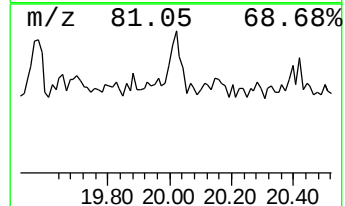
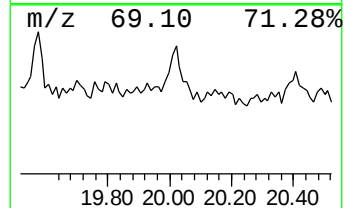
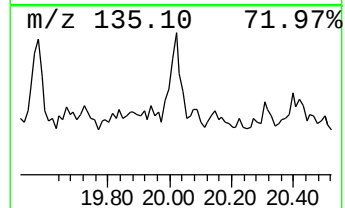
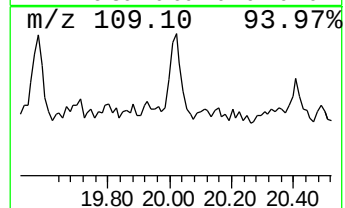
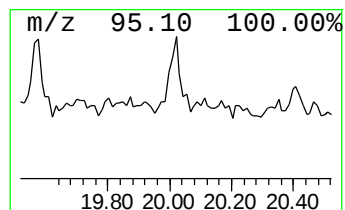
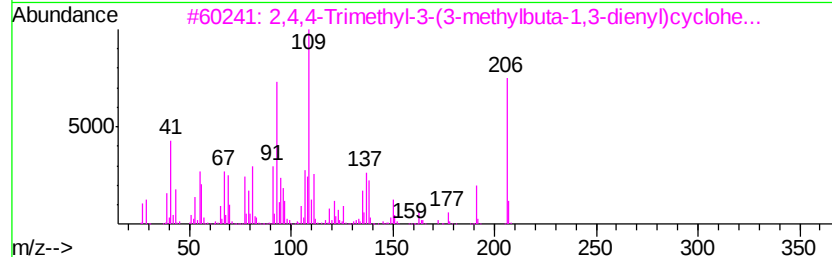
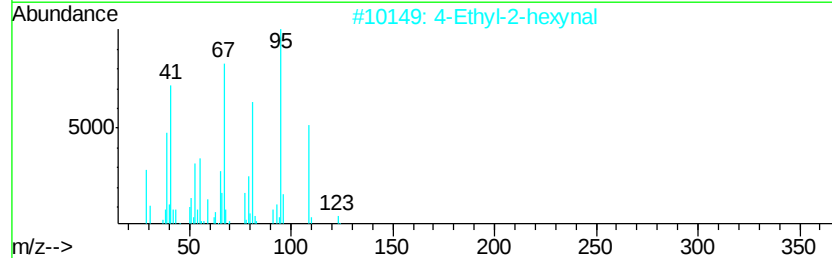
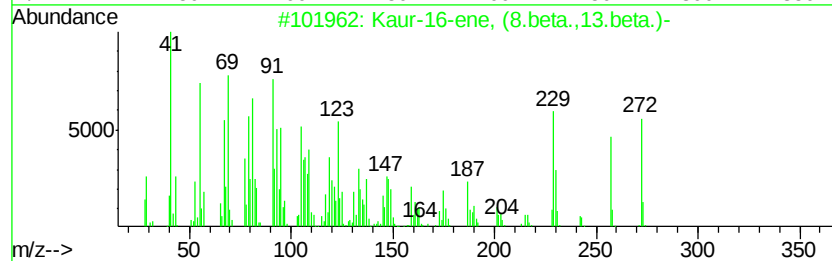
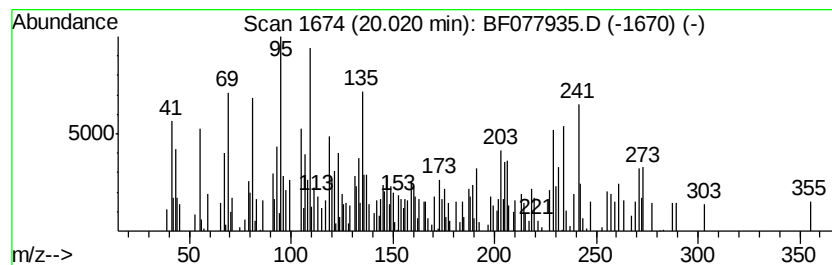
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown20.02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.02	6.41 ng	163619	Perylene-d12	17.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Kaur-16-ene, (8.beta.,13.beta.)-	272	C20H32	020070-61-5	38
2		4-Ethyl-2-hexynal	124	C8H12O	071932-97-3	27
3		2,4,4-Trimethyl-3-(3-methylbuta-...	206	C14H22O	097452-05-6	25
4		1-Oxaspiro[2.5]octane, 5,5-dimet...	206	C14H22O	1000195-92-1	25
5		Imidazole, 4-methyl-5-[3,3,3-tri...	234	C10H13F3N2O	1000129-15-8	20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077935.D
Acq On : 25 Mar 2015 4:35
Operator : TP/IZ
Sample : G1613-07
Misc :
ALS Vial : 26 Sample Multiplier: 1

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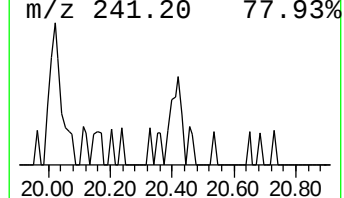
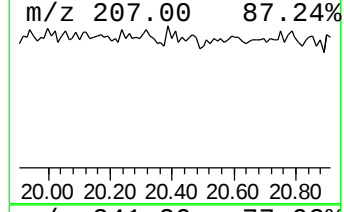
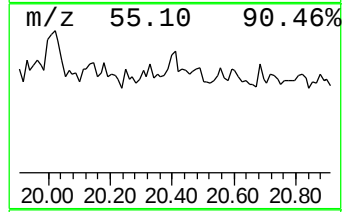
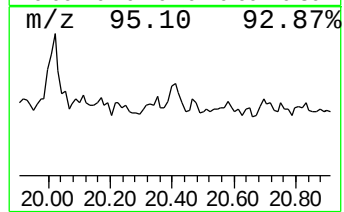
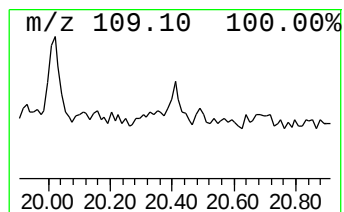
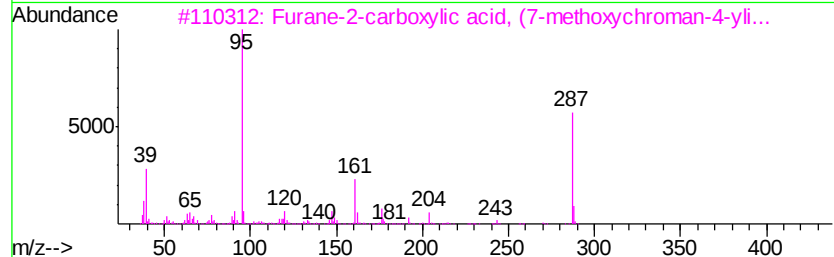
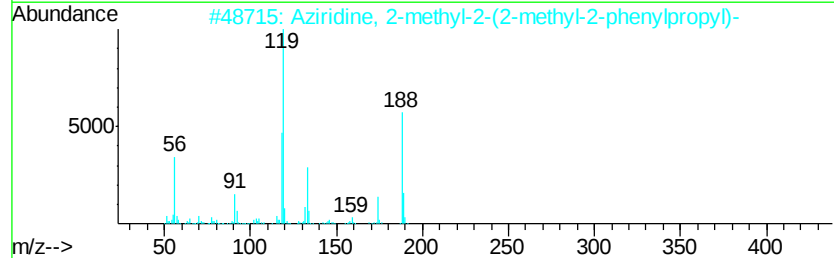
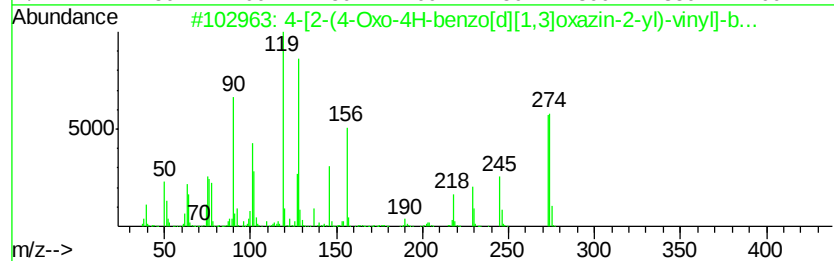
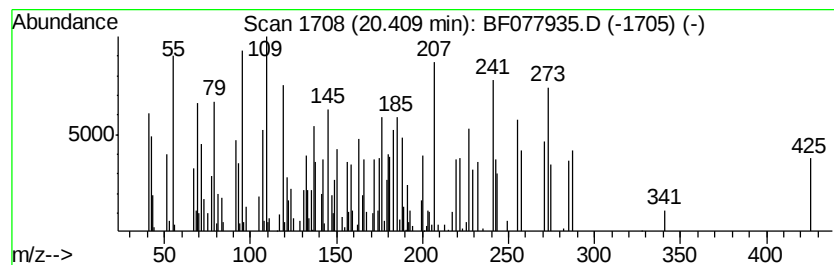
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown20.41 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.41	3.18 ng	81058	Perylene-d12	17.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-[2-(4-Oxo-4H-benzo[d][1,3]oxaz...	274	C17H10N2O2	1000297-20-9	11
2		Aziridine, 2-methyl-2-(2-methyl-...	189	C13H19N	1000293-28-4	11
3		Furane-2-carboxylic acid, (7-met...	287	C15H13NO5	332129-52-9	10
4		5.beta.-Cholest-23-ene, (Z)-	370	C27H46	014949-12-3	10
5		Cholan-24-oic acid, 3,12-dihydro...	406	C25H42O4	003245-38-3	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
 Data File : BF077935.D
 Acq On : 25 Mar 2015 4:35
 Operator : TP/IZ
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 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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
Cyclopentane, met...	1.05	9.8	ng	122185	1	7.12	250480	20.0
Propane, 2,2-dime...	1.28	2.7	ng	34311	1	7.12	250480	20.0
Butane, 2-methoxy...	1.56	4.5	ng	56126	1	7.12	250480	20.0
2-Pentanone, 4-hy...	4.86	8.7	ng	109130	1	7.12	250480	20.0
unknown6.83	6.83	62.2	ng	779421	1	7.12	250480	20.0
17-Pentatriacontene	15.88	3.9	ng	94323	5	15.97	489929	20.0
unknown17.28	17.28	2.7	ng	69865	6	17.72	510482	20.0
Naphthalene, 1,2,...	19.27	33.8	ng	862700	6	17.72	510482	20.0
unknown19.57	19.57	7.3	ng	185936	6	17.72	510482	20.0
unknown20.02	20.02	6.4	ng	163619	6	17.72	510482	20.0
unknown20.41	20.41	3.2	ng	81058	6	17.72	510482	20.0