LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

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	1.401	42	45	49	rVB	3753172	4267155	100.00%	15.907%
2	1.573	57	60	63	rVV	292189	354295	8.30%	1.321%
3	1.675	66	69	74	rVB3	19600	46498	1.09%	0.173%
4	1.767	75	77	88	rVB	122459	171398	4.02%	0.639%
5	2.315	121	125	128	rVB	25166	48805	1.14%	0.182%
6	2.475	136	139	147	rVB	153955	252595	5.92%	0.942%
7	2.978	179	183	188	rVB2	22419	48477	1.14%	0.181%
8	3.264	204	208	213	rBV	444939	849755	19.91%	3.168%
9	3.938	261	267	270	rBV	50253	88254	2.07%	0.329%
10	4.064	274	278	283	rVB	109224	172278	4.04%	0.642%
11	4.830	342	345	348	rBV	100885	164155	3.85%	0.612%
12	4.887	348	350	354	rVB2	29311	53754	1.26%	0.200%
13	5.207	375	378	383	rBV2	178922	279692	6.55%	1.043%
14	5.356	388	391	395	rVV	835003	1344023	31.50%	5.010%
15	5.413	395	396	399	rVB	316930	402944	9.44%	1.502%
16	5.619	405	414	417	rBV	50756	133936	3.14%	0.499%
17	5.676	417	419	422	rBV	484792	559794	13.12%	2.087%
18	5.962	442	444	447	rVB2	95684	124011	2.91%	0.462%
19	6.099	454	456	458	rBV	95447	105601	2.47%	0.394%
20	6.304	469	474	478	rVB2	36874	68841	1.61%	0.257%
21	6.544	492	495	497	rBV	75319	84260	1.97%	0.314%
22	6.636	501	503	505	rBV	100955	93942	2.20%	0.350%
23	6.670	505	506	508	rVV	86859	126564	2.97%	0.472%
24	6.704	508	509	511	rVV	282767	277240	6.50%	1.033%
25	6.739	511	512	516	rVB3	115966	113861	2.67%	0.424%
26	6.842	518	521	523	rVB	1017207	961989	22.54%	3.586%
27	6.910	523	527	531	rBV2	102454	200635	4.70%	0.748%
28	7.127	544	546	549	rBV	300939	285441	6.69%	1.064%
29	7.219	549	554	560	rVB2	117799	187089	4.38%	0.697%
30	7.310	560	562	564	rBV	996884	955347	22.39%	3.561%
31	7.585	569	586	588	rVB	494083	2693354	63.12%	10.040%
32	7.767	600	602	604	rVB	36611	43974	1.03%	0.164%
33	7.825	604	607	609	rBV	761725	807763	18.93%	3.011%
34	7.882	609	612	613	rVV	213665	330226	7.74%	1.231%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

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

35	7.985	613	621	623	rVB2	625943	2257536	52.90%	8.415%
36	8.190	635	639	641	rBV	339510	340592	7.98%	1.270%
37	8.225	641	642	646	rVB2	41688	60564	1.42%	0.226%
38	8.636	674	678	681	rBV3	26879	69652	1.63%	0.260%
39	8.705	681	684	686	rBV	367558	372258	8.72%	1.388%
40	8.990	703	709	711	rBV	60989	101796	2.39%	0.379%
41	9.082	715	717	719	rVB	46187	53788	1.26%	0.201%
42	9.185	722	726	729	rVV3	43954	88424	2.07%	0.330%
43	9.253	729	732	735	rVV3	41115	61629	1.44%	0.230%
44	9.368	739	742	744	rBV	40678	47158	1.11%	0.176%
45	9.642	761	766	769	rBV	68853	124322	2.91%	0.463%
46	9.768	774	777	780	rVV2	41978	117474	2.75%	0.438%
47	9.813	780	781	784	rVB2	39269	46706	1.09%	0.174%
48	9.905	786	789	792	rBV4	23772	61562	1.44%	0.229%
49	10.053	799	802	804	rVB	1717172	1609105	37.71%	5.998%
50	10.305	822	824	827	rVB	51482	49838	1.17%	0.186%
51	10.476	836	839	844	rVB2	45964	63413	1.49%	0.236%
52	10.876	871	874	876	rVB	379322	457093	10.71%	1.704%
53	11.848	957	959	962	rBV	902875	948444	22.23%	3.536%
54	12.705	1032	1034	1037	rBV	454028	454724	10.66%	1.695%
55	14.705	1206	1209	1211	rBV	1695378	1652148	38.72%	6.159%
56	15.997	1320	1322	1325	rBV	461920	493829	11.57%	1.841%
57	17.780	1475	1478	1481	rBV	413205	479719	11.24%	1.788%
58	20.226	1688	1692	1701	rVB	42260	116439	2.73%	0.434%

Sum of corrected areas: 26826159

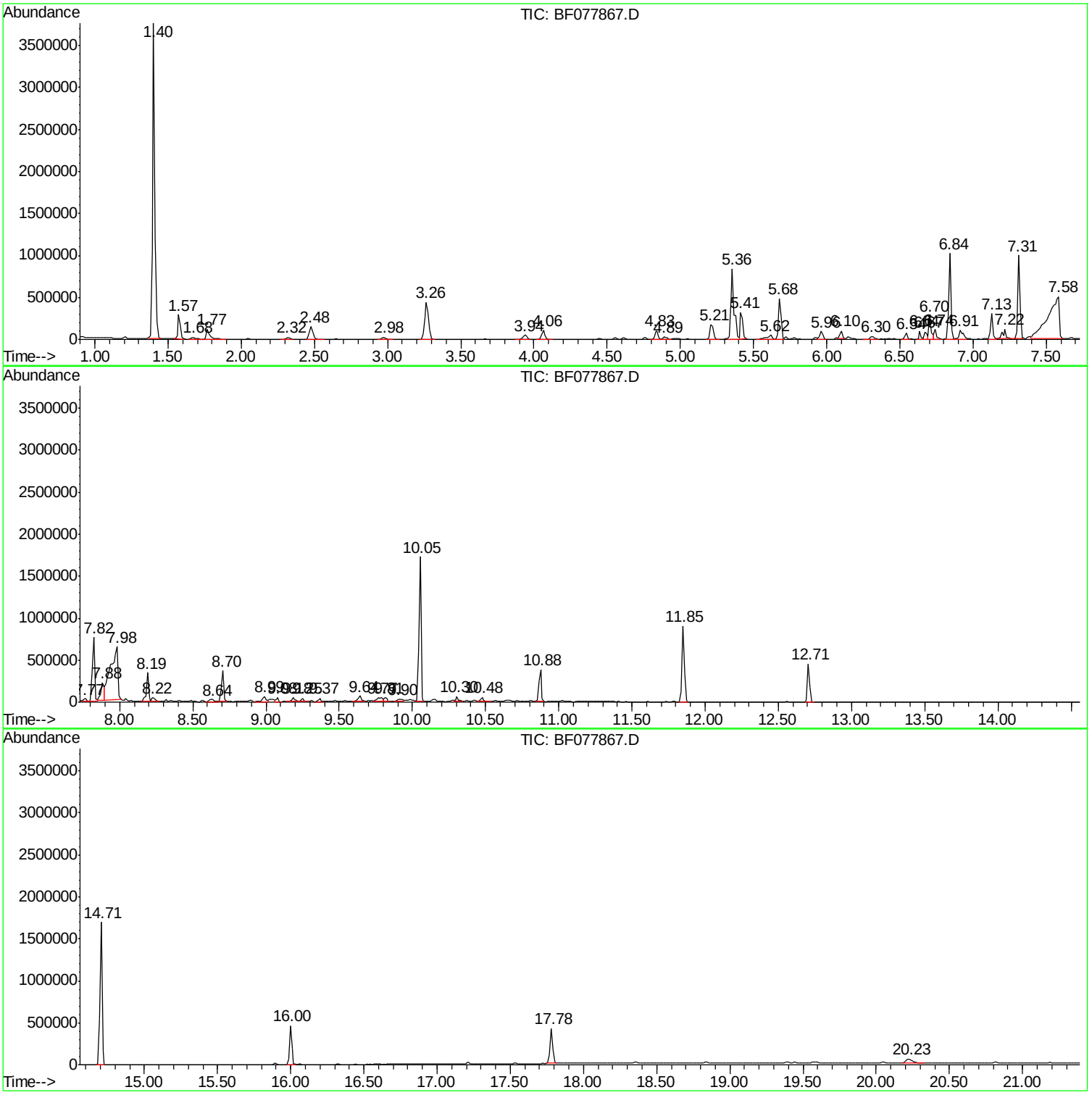
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

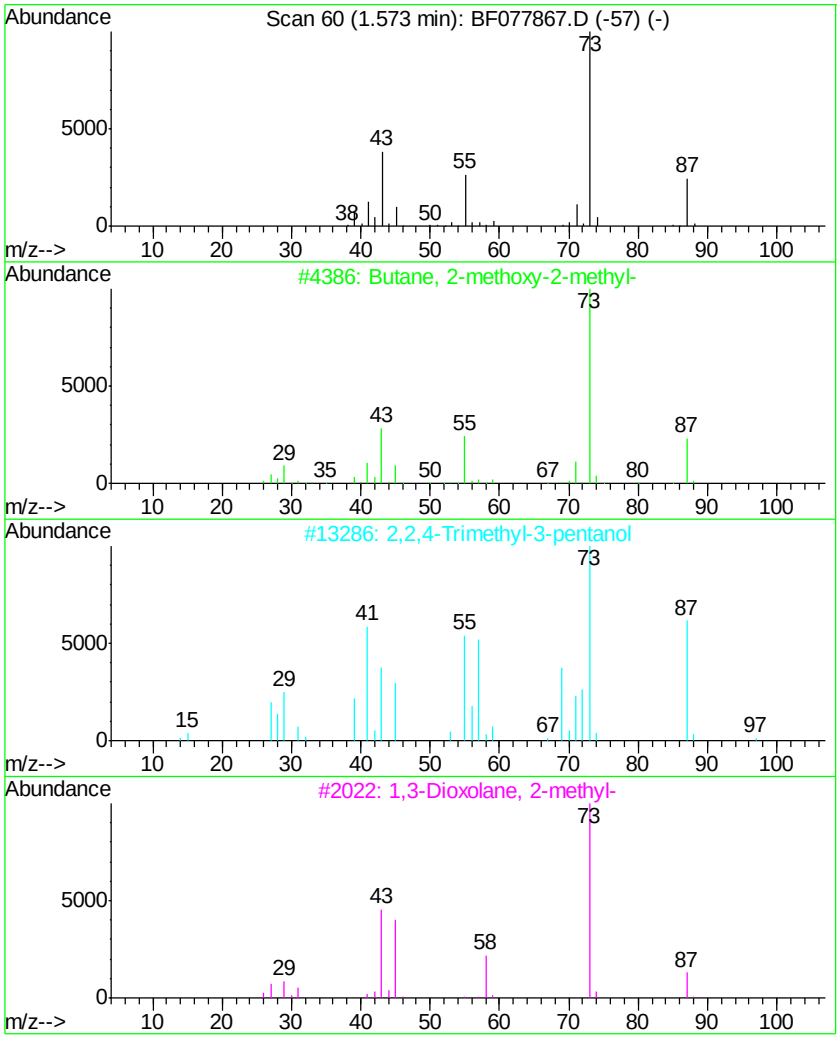
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

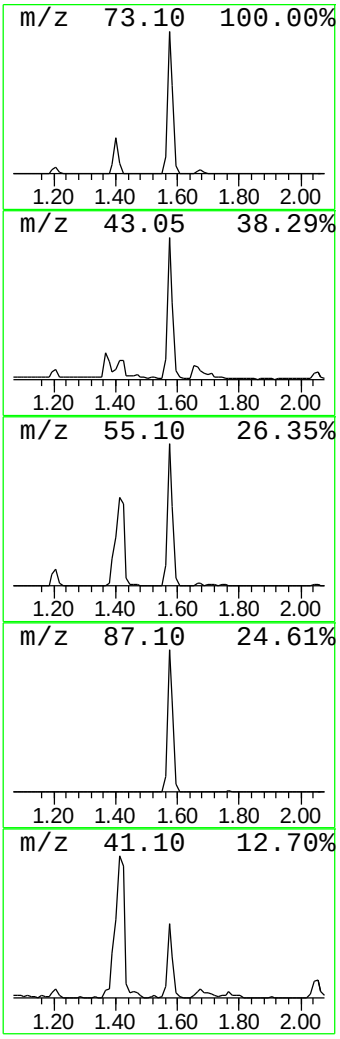
****

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.57	24.82 ng	354295	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

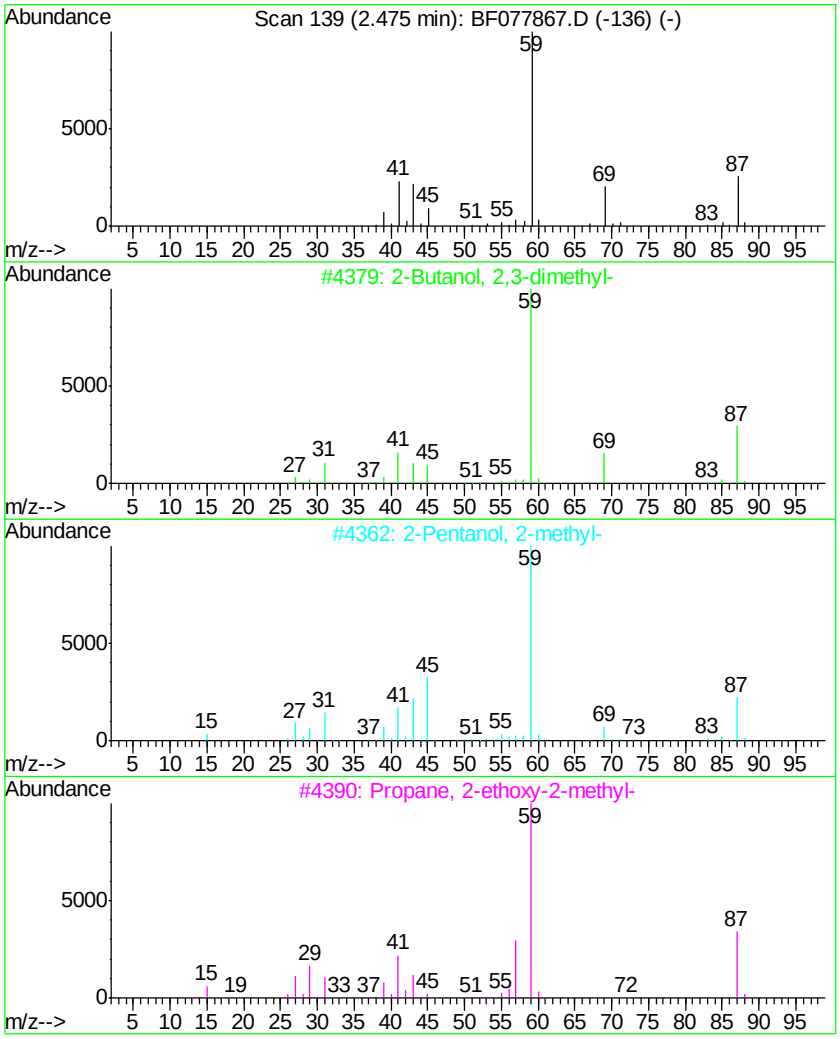
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

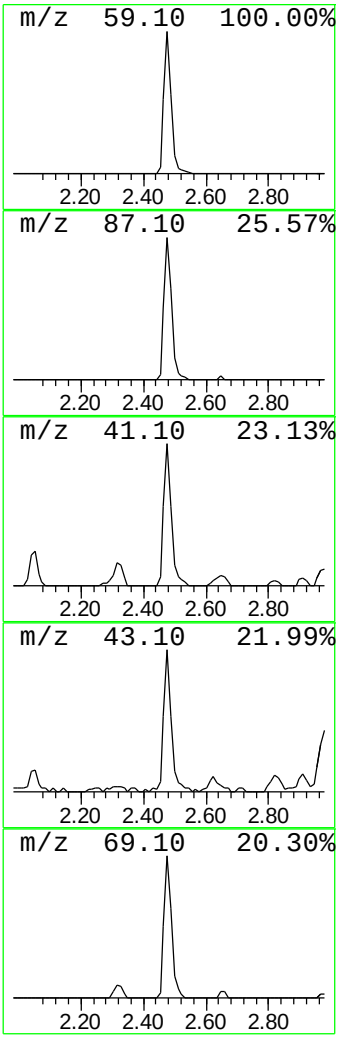
****

Peak Number 4 2-Butanol, 2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.48	17.70 ng	252595	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	56
3	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	50
4	2-Butanone, 3-ethoxy-3-methyl-	130	C7H14O2	036687-99-7	50
5	3-Heptanol	116	C7H16O	000589-82-2	40





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

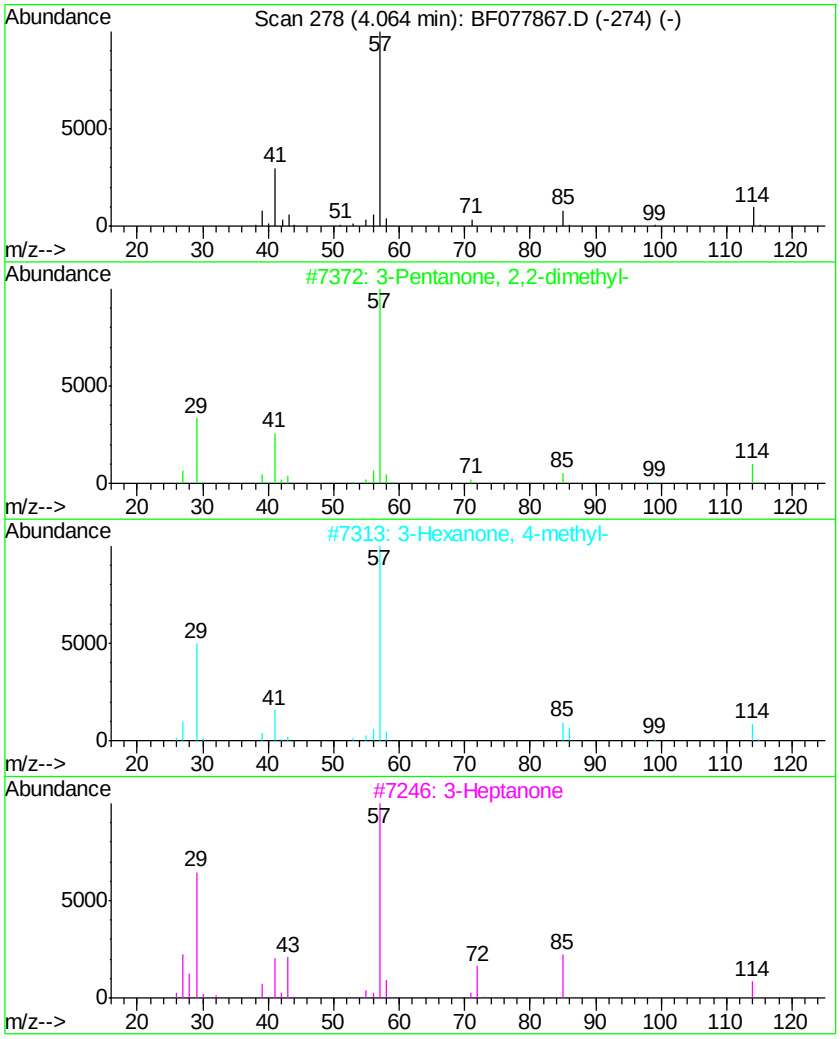
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

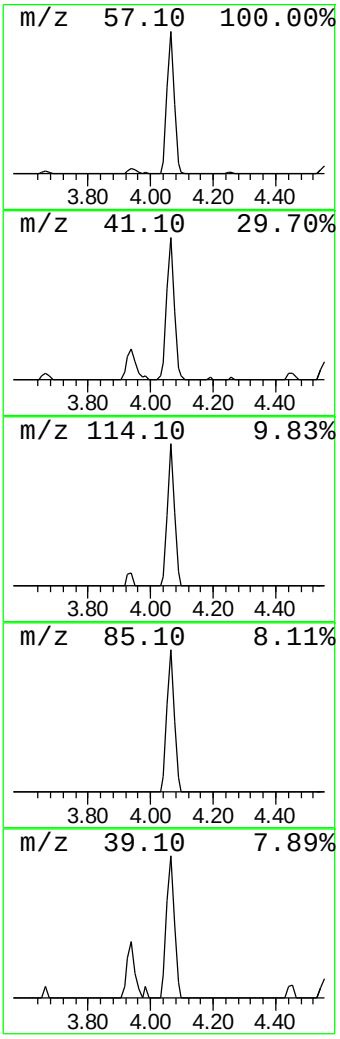
****

Peak Number 6 3-Pentanone, 2,2-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.06	12.07 ng	172278	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone, 2,2-dimethyl-	114	C7H14O	000564-04-5	87
2	3-Hexanone, 4-methyl-	114	C7H14O	017042-16-9	53
3	3-Heptanone	114	C7H14O	000106-35-4	50
4	2,2,5,5-Tetramethyl-3-hexanone	156	C10H20O	000868-91-7	40
5	3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

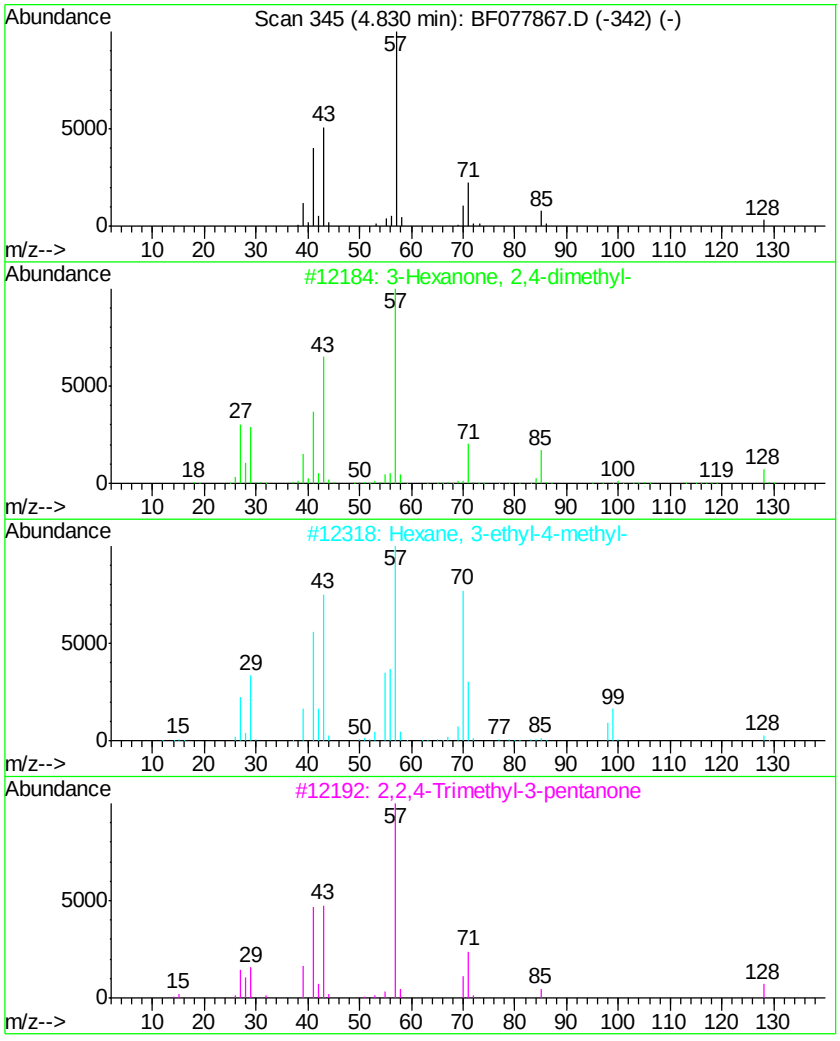
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

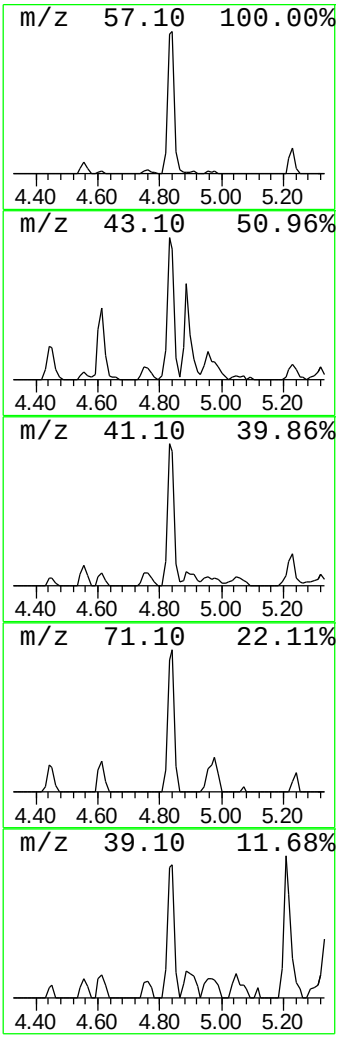
****

Peak Number 7 3-Hexanone, 2,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	11.50 ng	164155	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone, 2,4-dimethyl-	128	C8H16O	018641-70-8	53
2	Hexane, 3-ethyl-4-methyl-	128	C9H20	003074-77-9	53
3	2,2,4-Trimethyl-3-pentanone	128	C8H16O	005857-36-3	50
4	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	40
5	Octane, 2,2-dimethyl-	142	C10H22	015869-87-1	40





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

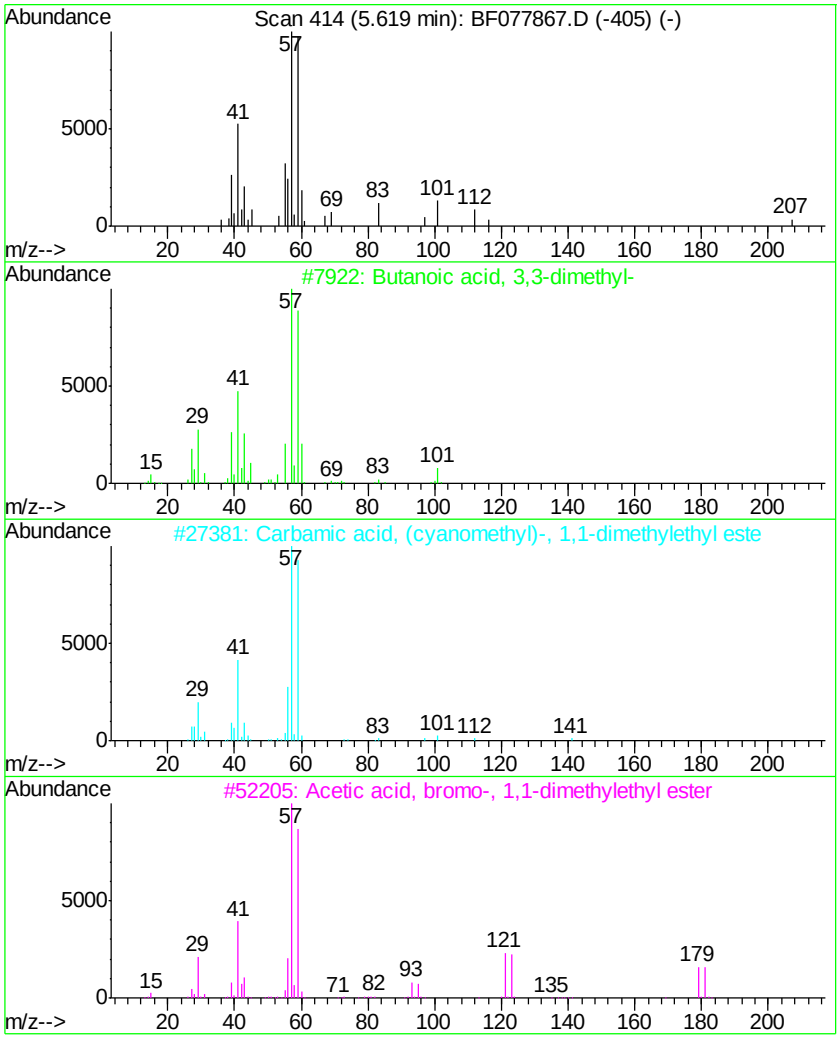
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

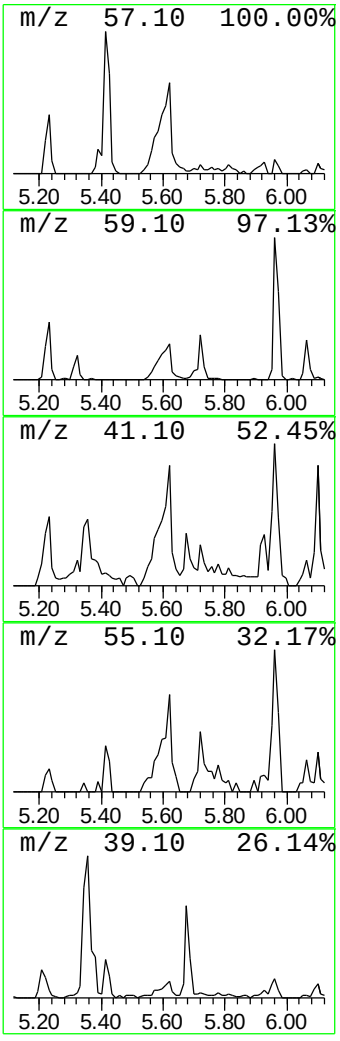
****

Peak Number 10 Butanoic acid, 3,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.62	9.38 ng	133936	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3,3-dimethyl-	116	C6H12O2	001070-83-3	53
2	Carbamic acid, (cyanomethyl)-, 1...	156	C7H12N2O2	085363-04-8	50
3	Acetic acid, bromo-, 1,1-dimethy...	194	C6H11BrO2	005292-43-3	42
4	Acetic acid, cyano(hydroxyimino)...	170	C7H10N2O3	076101-98-9	42
5	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

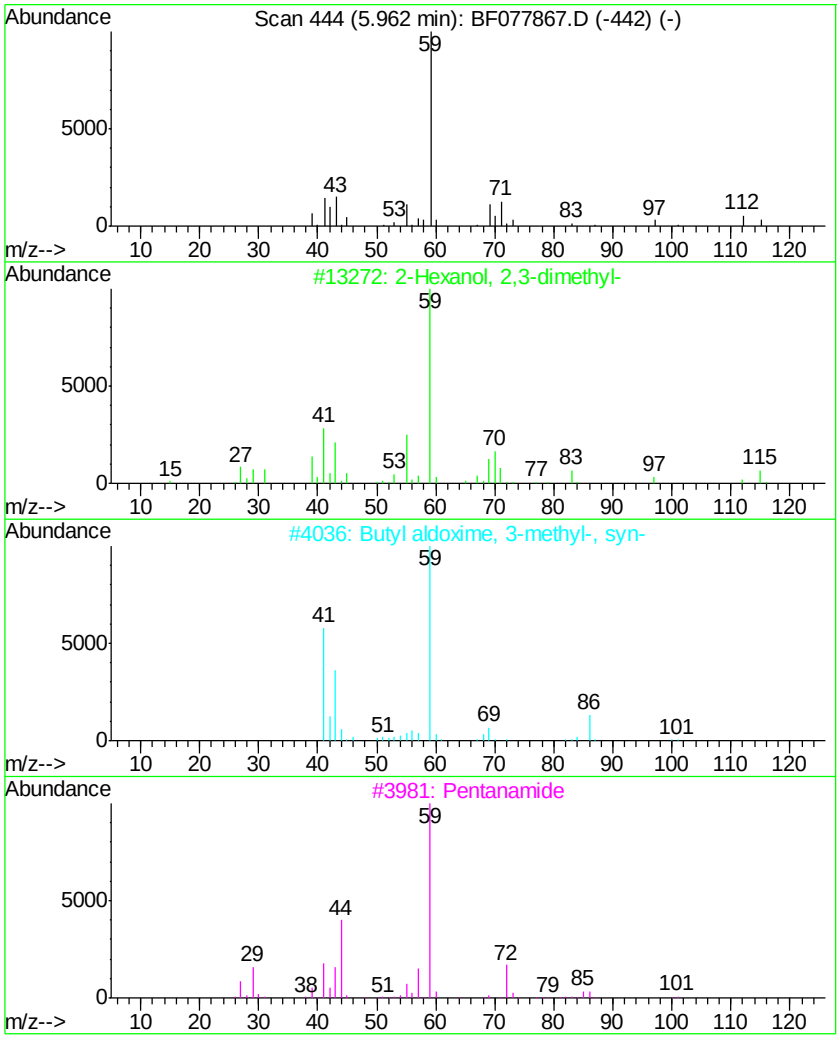
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

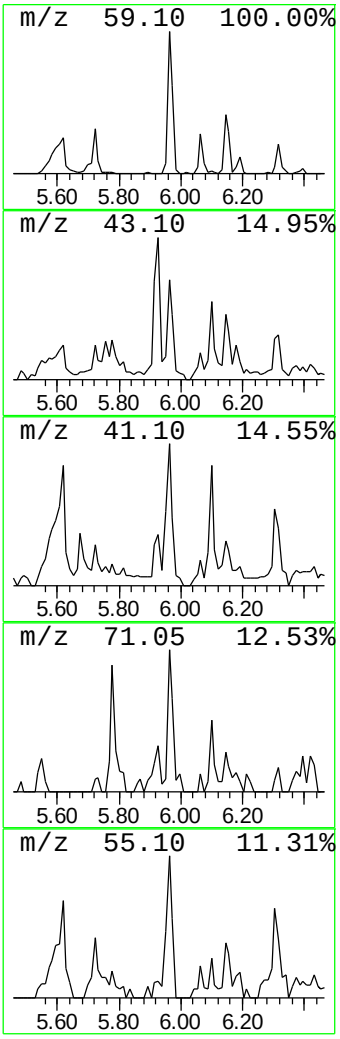
****

Peak Number 12 2-Hexanol, 2,3-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.96	8.69 ng	124011	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	72
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	59
3	Pentanamide	101	C5H11NO	000626-97-1	59
4	Butyl aldoxime, 3-methyl-, anti	101	C5H11NO	005775-74-6	50
5	2-Hydroxy-2-methylundec-5-en-3-one	198	C12H22O2	1000191-46-2	39





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

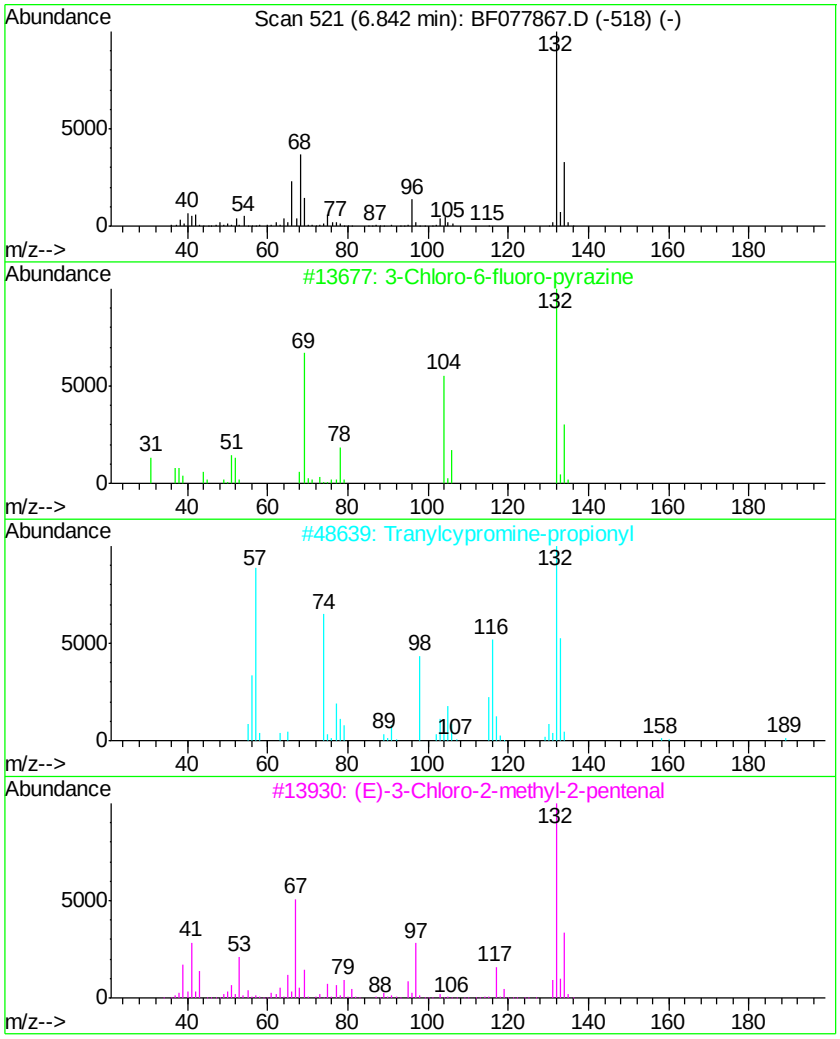
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

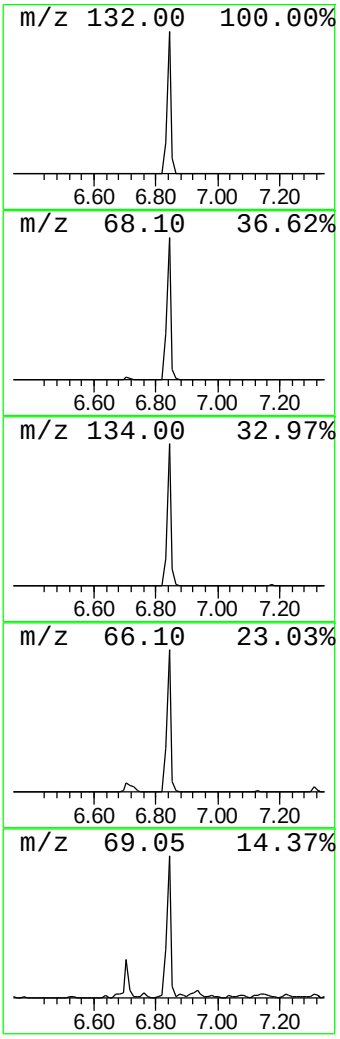
****

Peak Number 13 unknown6.84 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.84	67.40 ng	961989	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	12
5	Xenon	132	Xe	007440-63-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

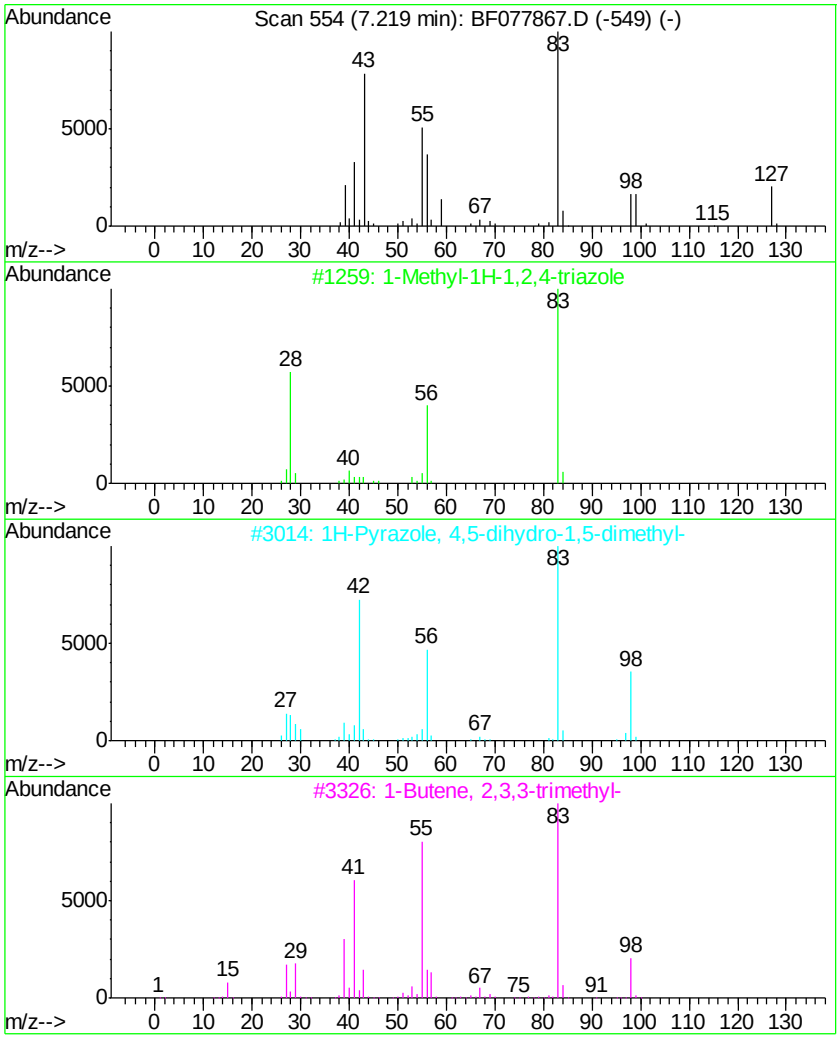
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

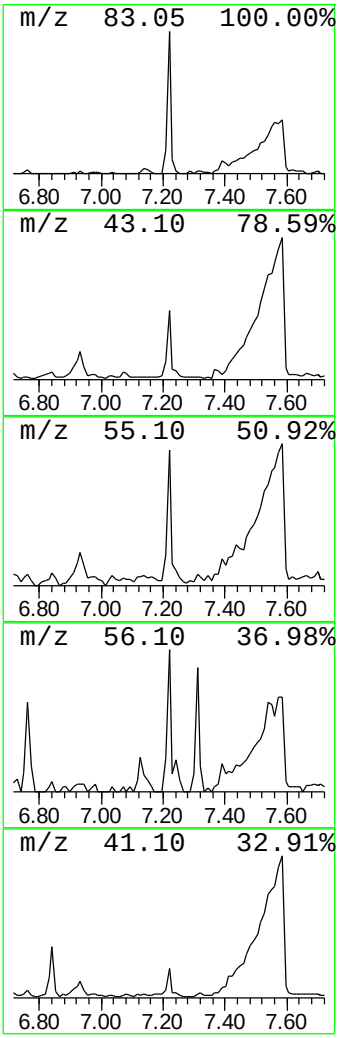
****

Peak Number 15 unknown7.22 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	13.11 ng	187089	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	47
2	1H-Pyrazole, 4,5-dihydro-1,5-dim...	98	C5H10N2	005775-96-2	47
3	1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	47
4	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	43
5	2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

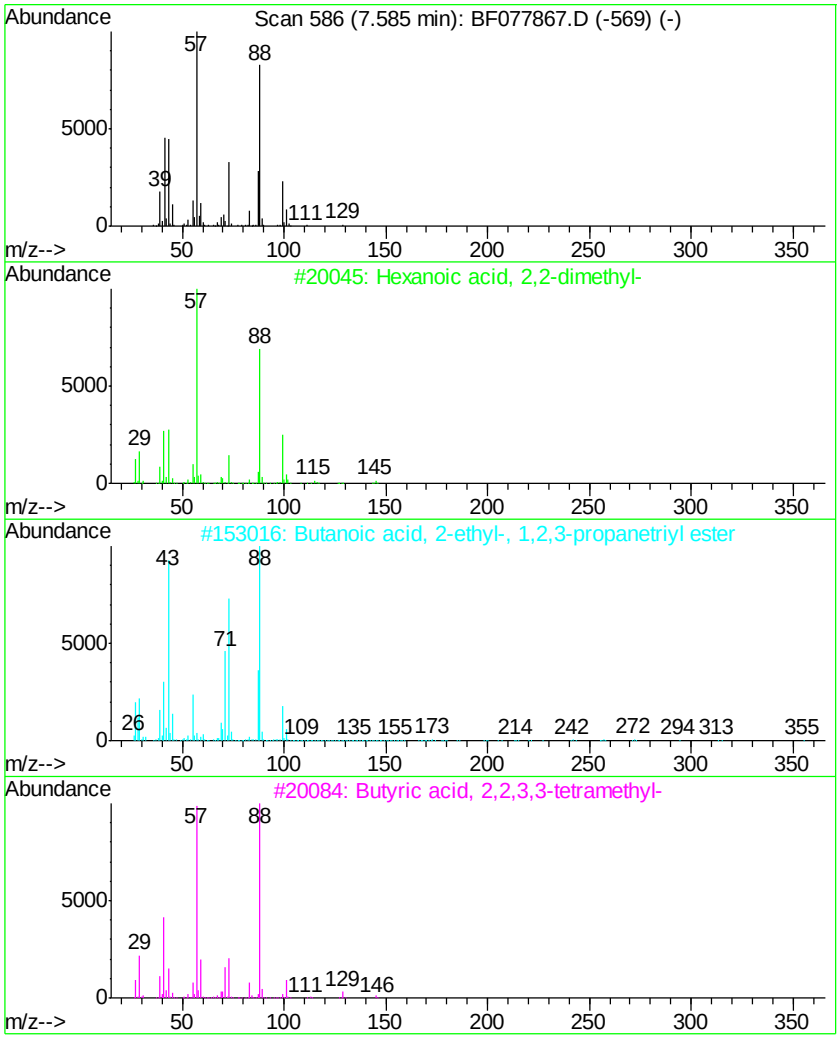
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

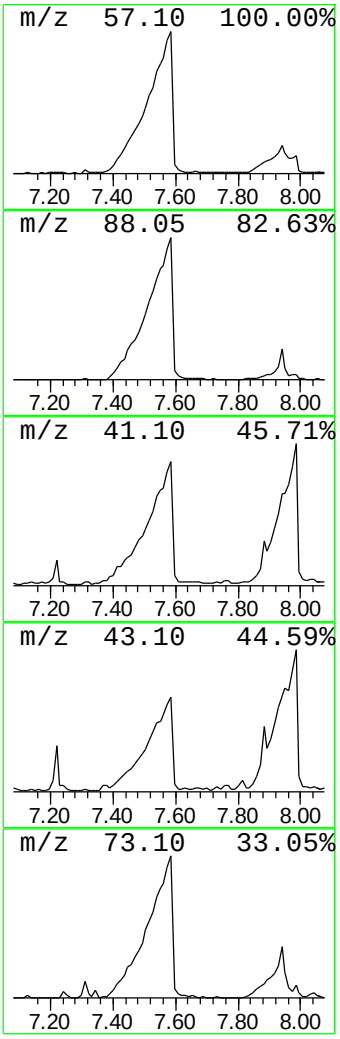
****

Peak Number 17 Hexanoic acid, 2,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	188.72 ng	2693350	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2,2-dimethyl-	144	C8H16O2	000813-72-9	59
2	Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	45
3	Butyric acid, 2,2,3,3-tetramethyl-	144	C8H16O2	030407-41-1	42
4	Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	40
5	Methyl propionate	88	C4H8O2	000554-12-1	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

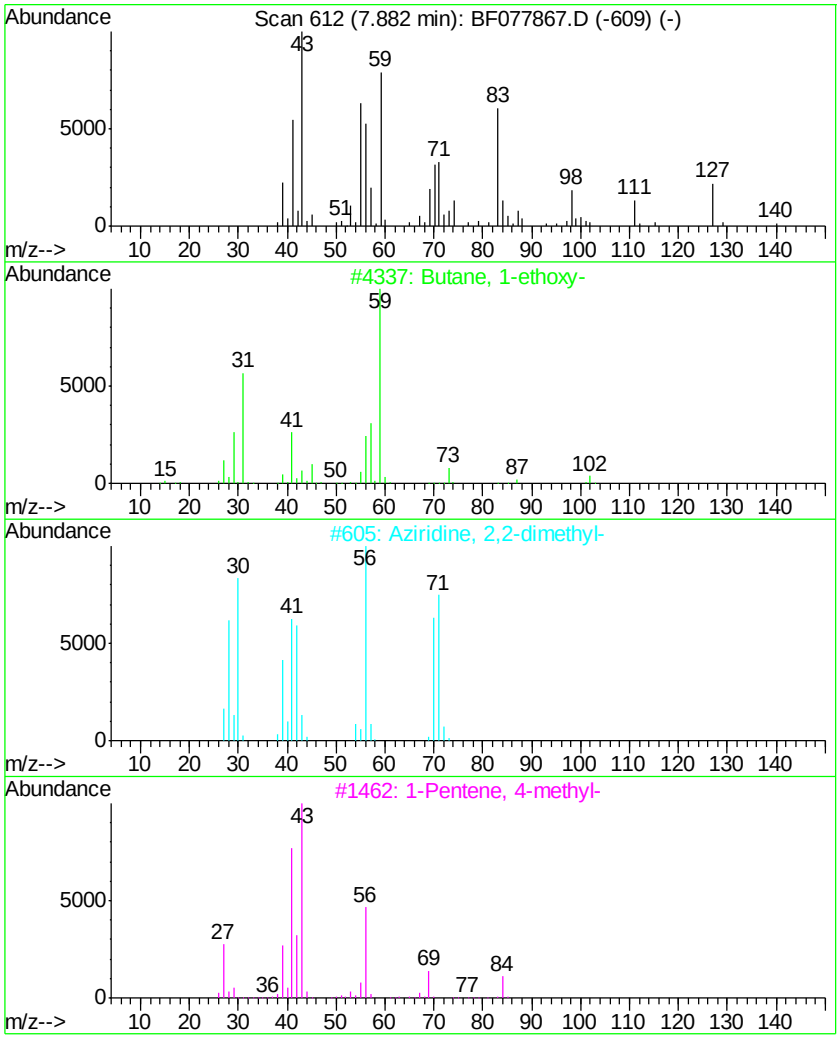
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

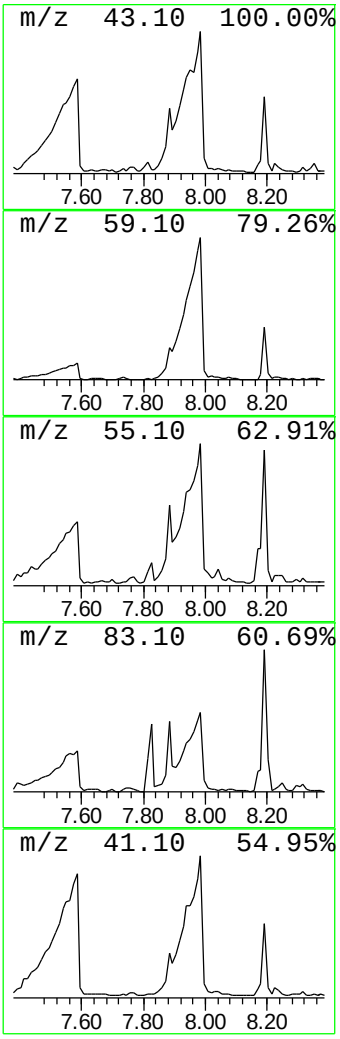
****

Peak Number 18 unknown7.88 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	23.14 ng	330226	1,4-Dichlorobenzene-d4	7.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	22
2	Aziridine, 2,2-dimethyl-	71	C4H9N	002658-24-4	14
3	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	11
4	2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	11
5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	11





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

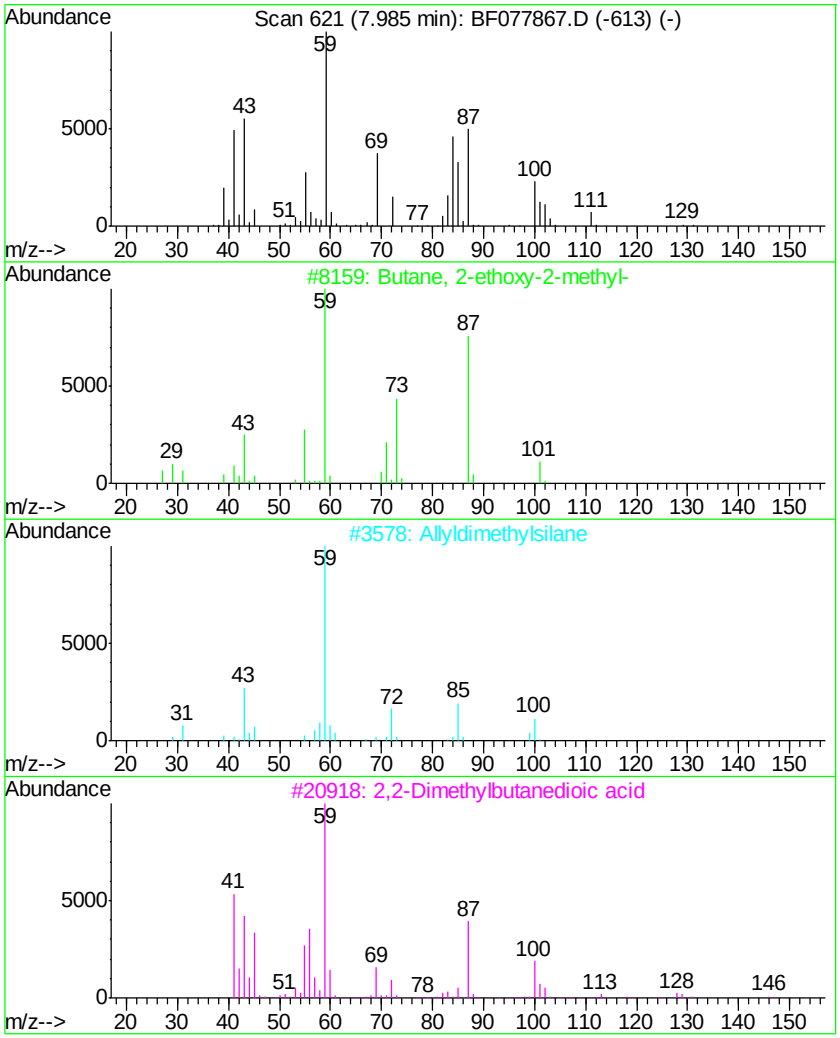
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

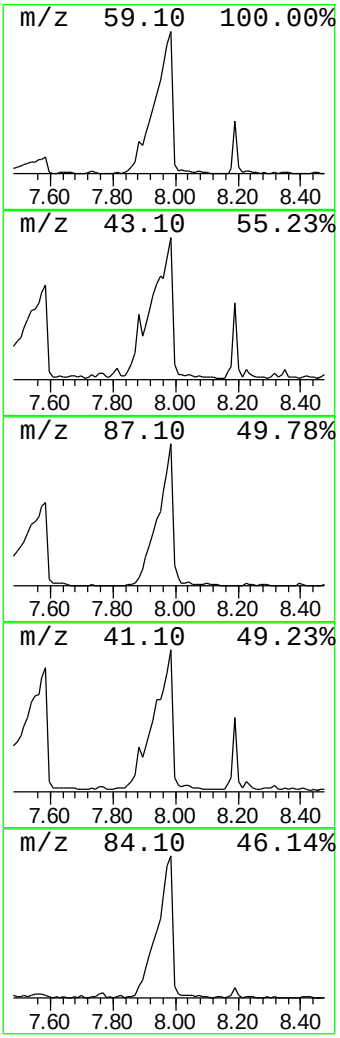
****

Peak Number 19 unknown7.98 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.98	121.29 ng	2257540	Naphthalene-d8	8.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-ethoxy-2-methyl-	116	C7H16O	000919-94-8	37
2	Allyldimethylsilane	100	C5H12Si	003937-30-2	37
3	2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	36
4	1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	27
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	23





Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

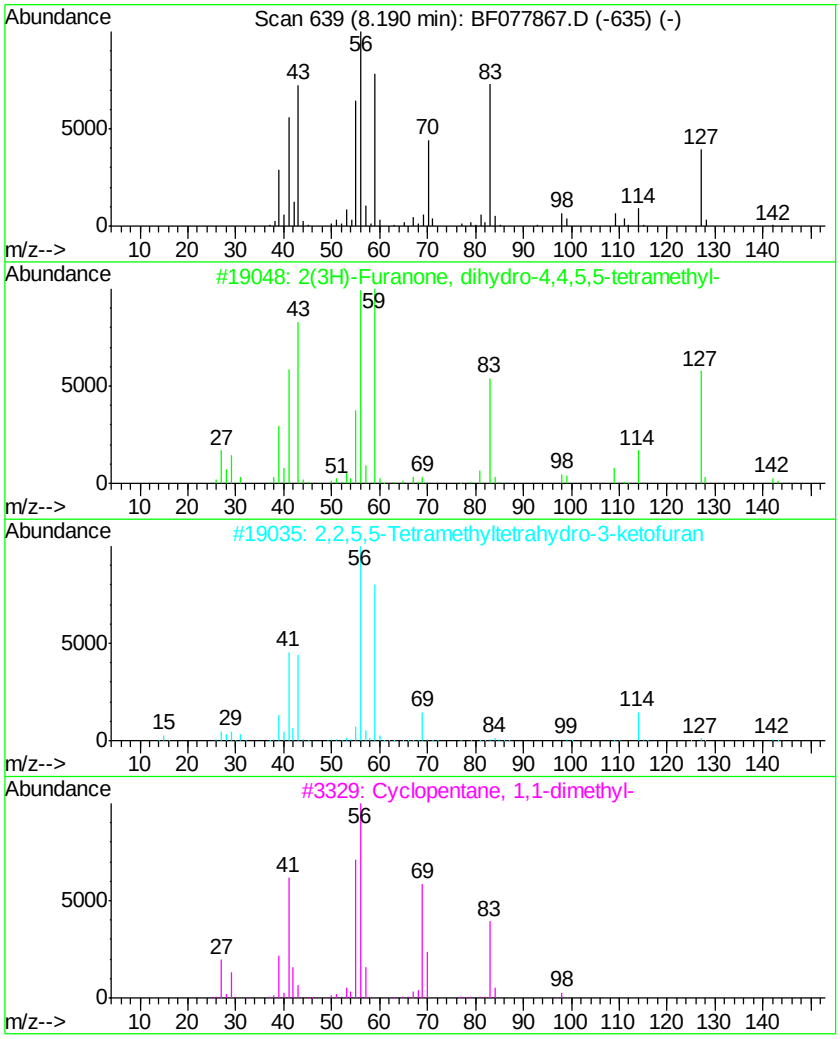
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

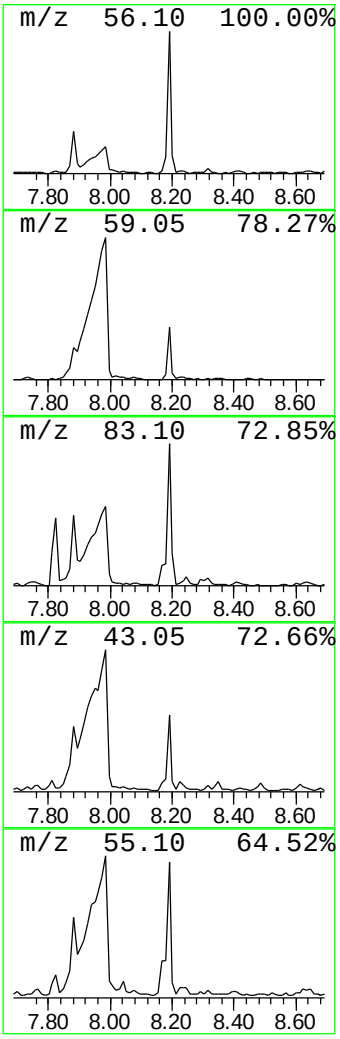
****

Peak Number 20 unknown8.19 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.19	18.30 ng	340592	Naphthalene-d8	8.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Furanone, dihydro-4,4,5,5-...	142	C8H14O2	016466-24-3	43
2	2,2,5,5-Tetramethyltetrahydro-3-...	142	C8H14O2	005455-94-7	32
3	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	30
4	2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	27
5	Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032015\
Data File : BF077867.D
Acq On : 20 Mar 2015 17:04
Operator : TP/IZ
Sample : G1604-01
Misc : W/DOD
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : Z2-0121

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	1.57	24.8	ng	354295	1	7.13	285441	20.0
2-Butanol, 2,3-di...	2.48	17.7	ng	252595	1	7.13	285441	20.0
3-Pentanone, 2,2-...	4.06	12.1	ng	172278	1	7.13	285441	20.0
3-Hexanone, 2,4-d...	4.83	11.5	ng	164155	1	7.13	285441	20.0
Butanoic acid, 3,...	5.62	9.4	ng	133936	1	7.13	285441	20.0
2-Hexanol, 2,3-di...	5.96	8.7	ng	124011	1	7.13	285441	20.0
unknown6.84	6.84	67.4	ng	961989	1	7.13	285441	20.0
unknown7.22	7.22	13.1	ng	187089	1	7.13	285441	20.0
Hexanoic acid, 2,...	7.58	188.7	ng	2693350	1	7.13	285441	20.0
unknown7.88	7.88	23.1	ng	330226	1	7.13	285441	20.0
unknown7.98	7.98	121.3	ng	2257540	2	8.70	372258	20.0
unknown8.19	8.19	18.3	ng	340592	2	8.70	372258	20.0