LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016666.D
Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : RE114D1-20210916

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BN016666.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.161	4	9	11	rBV	711997	1128602	100.00%	27.713%
2	3.364	29	31	47	rVB	75820	232441	20.60%	5.708%
3	3.816	77	80	96	rVB	8074	34586	3.06%	0.849%
4	4.010	98	101	119	rVB	6529	25869	2.29%	0.635%
5	4.821	185	189	200	rVB	51363	118327	10.48%	2.906%
6	5.264	234	237	244	rVB	5309	13874	1.23%	0.341%
7	6.296	345	349	354	rBV	14599	31815	2.82%	0.781%
8	6.637	383	386	401	rVV	6313	15832	1.40%	0.389%
9	6.850	403	409	418	rVB3	5693	22287	1.97%	0.547%
10	7.080	430	434	440	rVB	5865	11708	1.04%	0.287%
11	7.264	449	454	460	rVB3	7782	21739	1.93%	0.534%
12	7.633	489	494	499	rVB2	52278	108847	9.64%	2.673%
13	7.864	515	519	526	rBV	19253	39159	3.47%	0.962%
14	8.168	549	552	559	rVV	11504	29002	2.57%	0.712%
15	8.306	563	567	569	rVB	16192	20753	1.84%	0.510%
16	8.785	601	605	609	rBV	24286	48729	4.32%	1.197%
17	9.114	626	631	635	rVB2	10205	25677	2.28%	0.631%
18	10.191	713	716	729	rBV	177265	319338	28.30%	7.841%
19	10.407	729	733	736	rBV	106408	182621	16.18%	4.484%
20	12.003	918	931	942	rBV	46856	76046	6.74%	1.867%
21	12.886	1071	1074	1078	rBV	105940	171125	15.16%	4.202%
22	14.269	1180	1183	1186	rBV	135564	203341	18.02%	4.993%
23	14.434	1193	1196	1198	rBV	11146	17062	1.51%	0.419%
24	15.715	1294	1297	1299	rBV	15161	21438	1.90%	0.526%
25	15.766	1299	1301	1305	rVB2	12809	20604	1.83%	0.506%
26	17.018	1400	1403	1406	rBV	111450	167746	14.86%	4.119%
27	19.063	1688	1696	1714	rBV	137801	188535	16.71%	4.630%
28	19.679	1813	1820	1840	rBV	135571	173199	15.35%	4.253%
29	20.983	1975	1978	1991	rBV2	6476	20711	1.84%	0.509%
30	21.164	1993	1995	1999	rBV	81270	121267	10.74%	2.978%
31	21.238	1999	2002	2011	rVB2	124095	202138	17.91%	4.964%
32	22.311	2100	2103	2108	rBV	10938	22226	1.97%	0.546%
33	23.485	2415	2427	2463	rVB	106817	209510	18.56%	5.145%
34	24.087	2593	2603	2620	rBV	6337	12699	1.13%	0.312%
35	24.854	2817	2827	2850	rVB	5896	13617	1.21%	0.334%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Sum of corrected areas: 4072470

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Acq On : 02 Oct 2021 11:08
Operator : CG/JU
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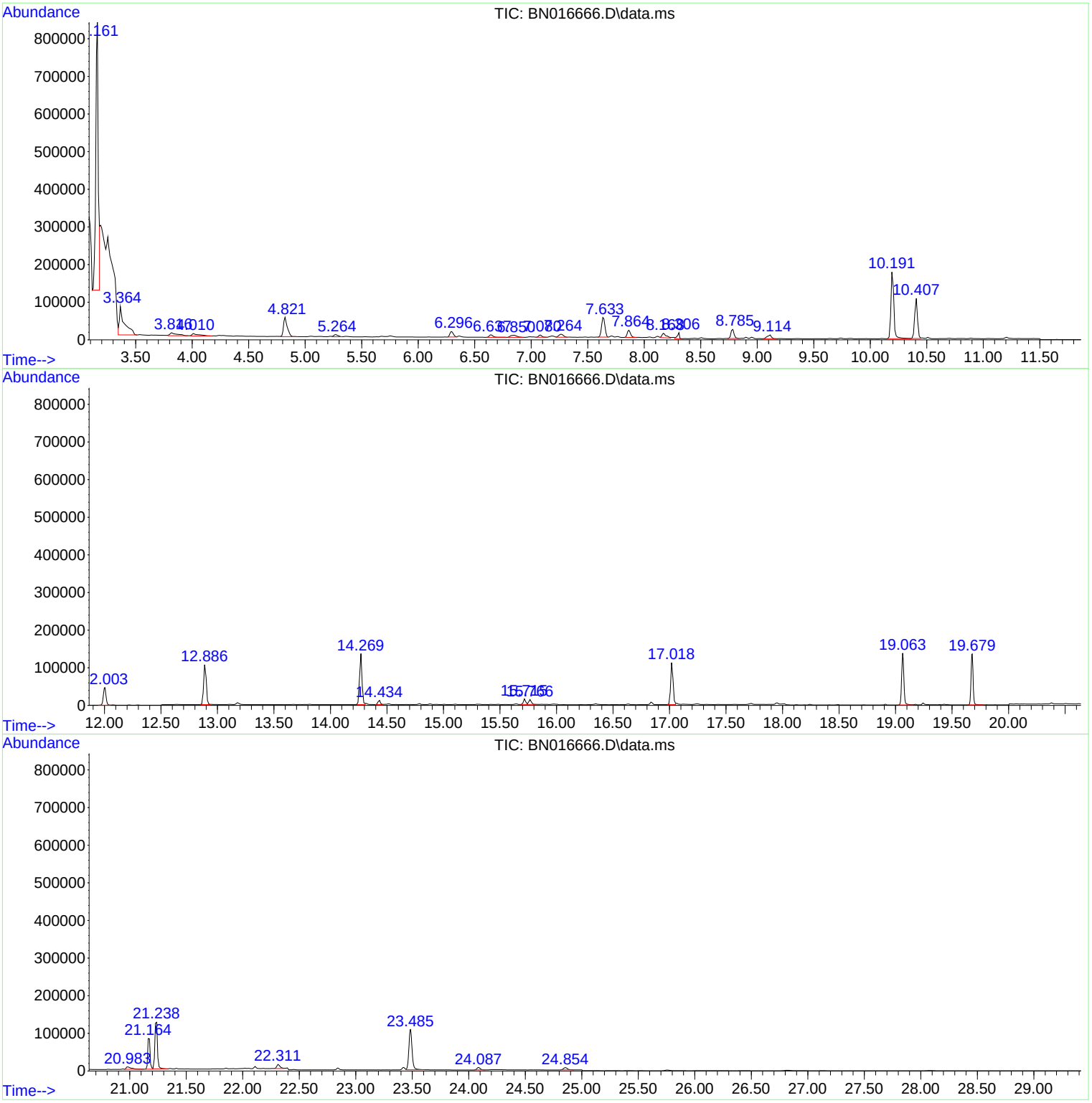
Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

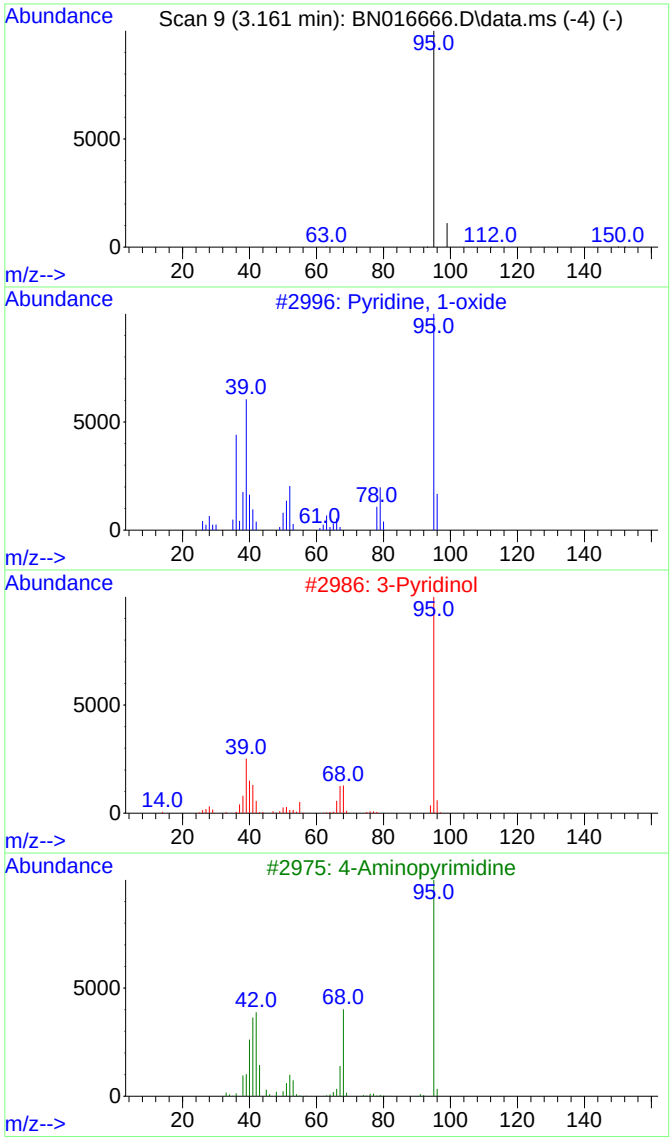
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

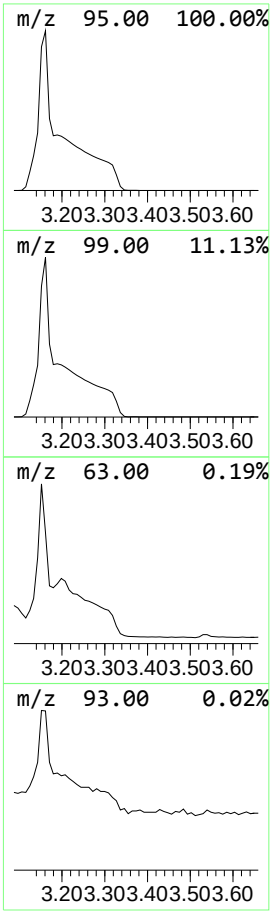
****

Peak Number 1 Pyridine, 1-oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.161	4.15 ng	1128600	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 1-oxide	95	C5H5NO	000694-59-7	2
2	3-Pyridinol	95	C5H5NO	000109-00-2	2
3	4-Aminopyrimidine	95	C4H5N3	000591-54-8	2
4	4-Pyridinol	95	C5H5NO	000626-64-2	2
5	2-(Aminomethylene)aminoacrylontrile	95	C4H5N3	167320-31-2	2





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016666.D
Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

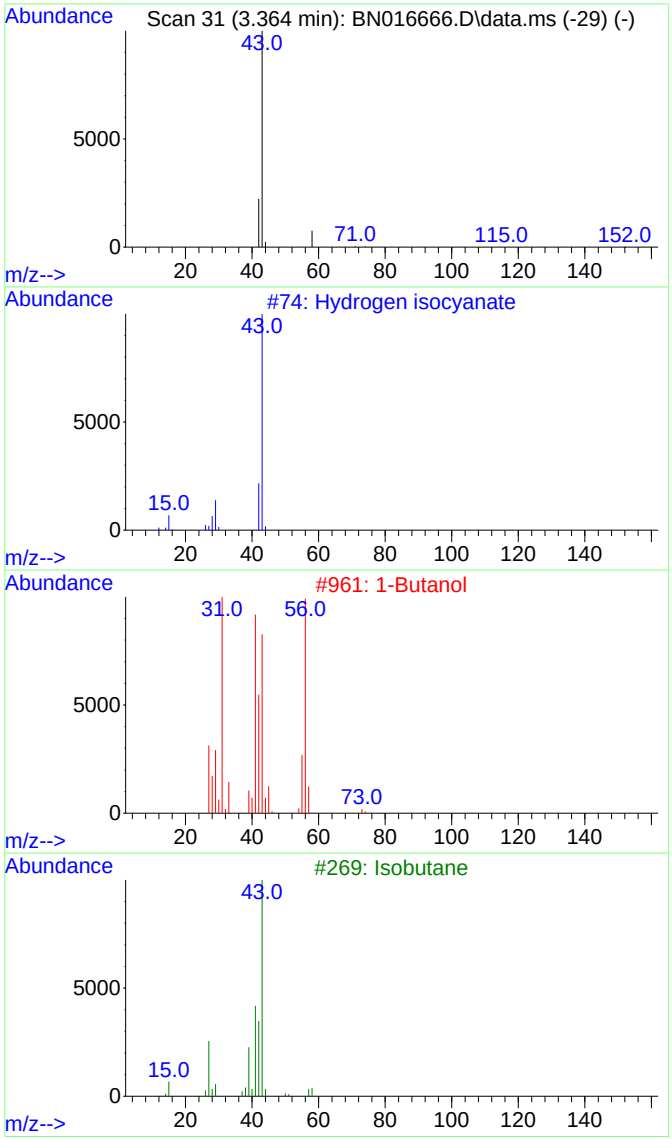
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

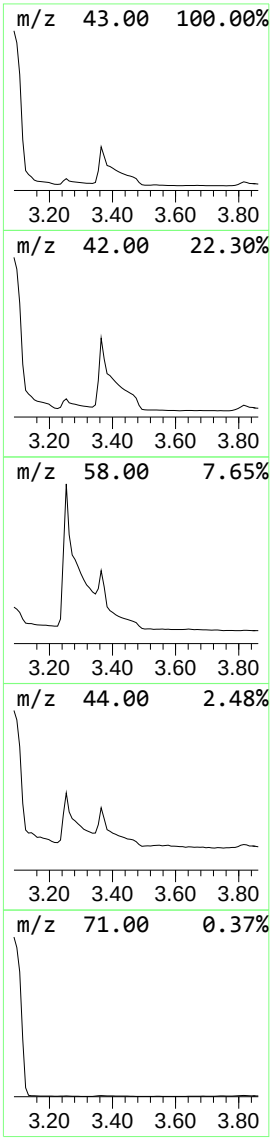
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Hydrogen isocyanate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.364	0.85 ng	232441	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrogen isocyanate	43	CHNO	000075-13-8	4
2	1-Butanol	74	C4H10O	000071-36-3	3
3	Isobutane	58	C4H10	000075-28-5	3
4	Hydrogen azide	43	HN3	007782-79-8	2
5	Butane	58	C4H10	000106-97-8	2





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

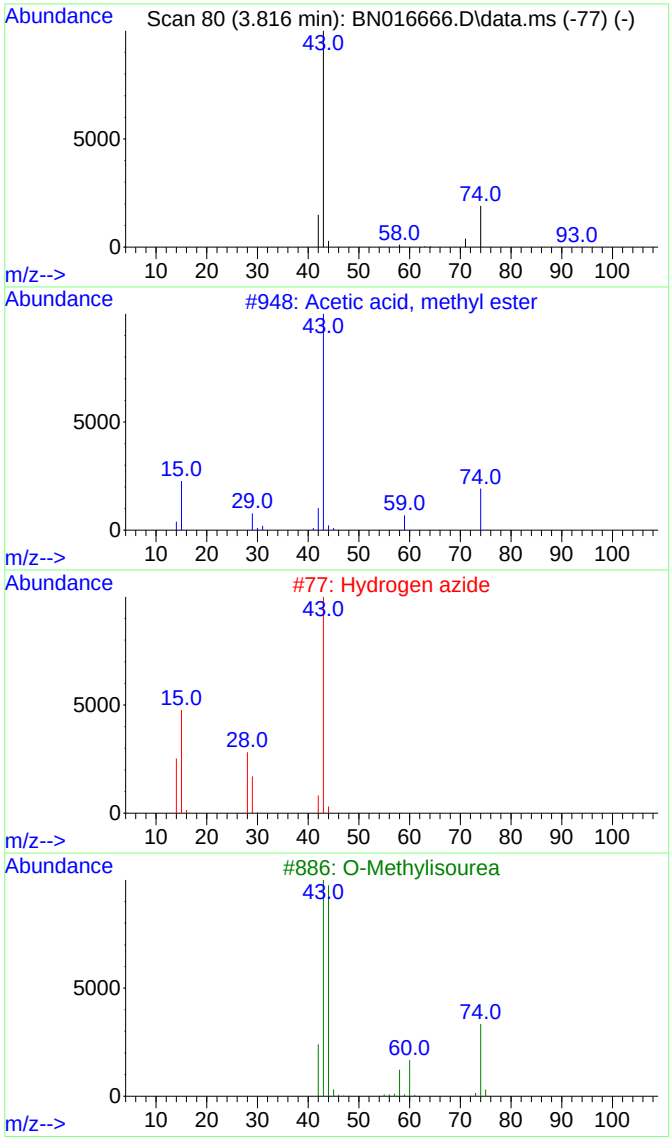
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

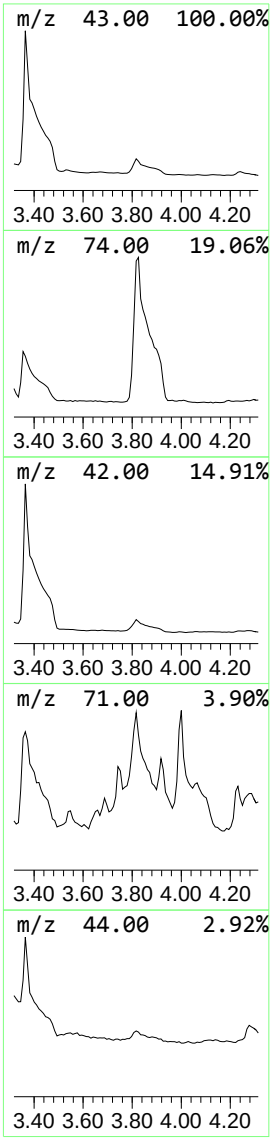
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Acetic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.816	0.13 ng	34586	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, methyl ester	74	C3H6O2	000079-20-9	9
2	Hydrogen azide	43	HN3	007782-79-8	4
3	O-Methylisourea	74	C2H6N2O	002440-60-0	4
4	Hydrogen isocyanate	43	CHNO	000075-13-8	3
5	1-Ethyl-1-methylhydrazine	74	C3H10N2	004986-48-5	3





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : RE114D1-20210916

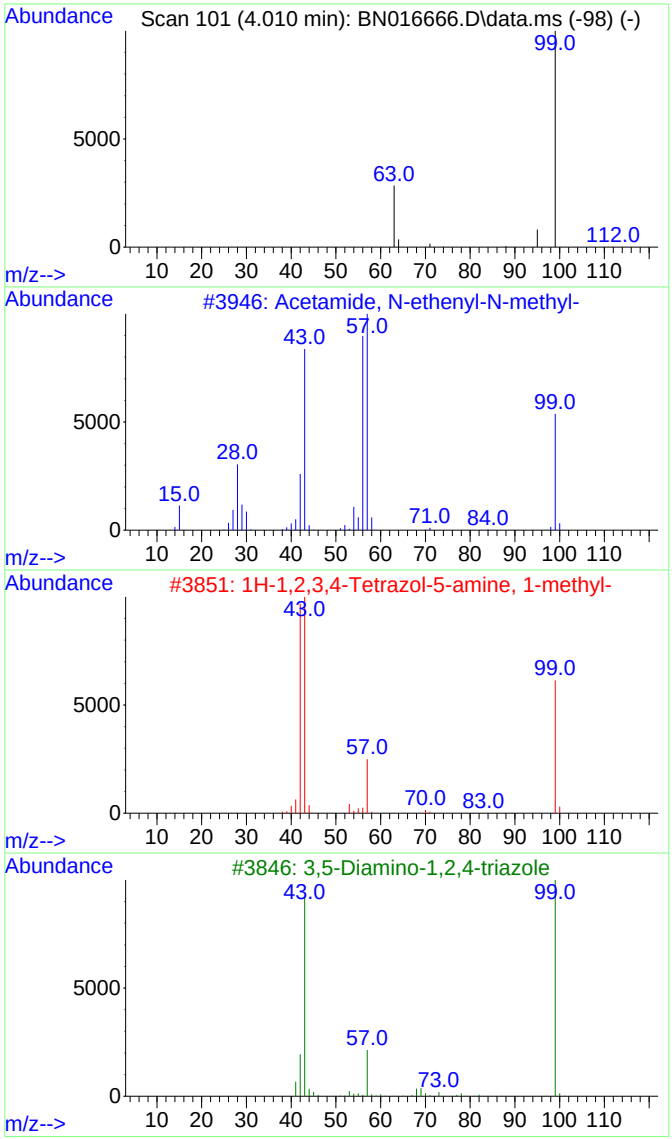
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

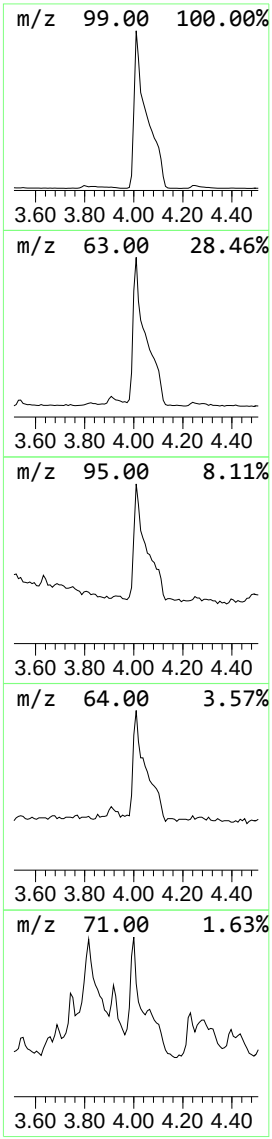
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Acetamide, N-ethenyl-N-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.010	0.10 ng	25869	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-ethenyl-N-methyl-	99	C5H9NO	003195-78-6	3
2	1H-1,2,3,4-Tetrazol-5-amine, 1-m...	99	C2H5N5	005422-44-6	3
3	3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	2
4	Isothiazole, 3-methyl-	99	C4H5NS	000693-92-5	2
5	Isothiazole, 4-methyl-	99	C4H5NS	000693-90-3	2





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
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Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

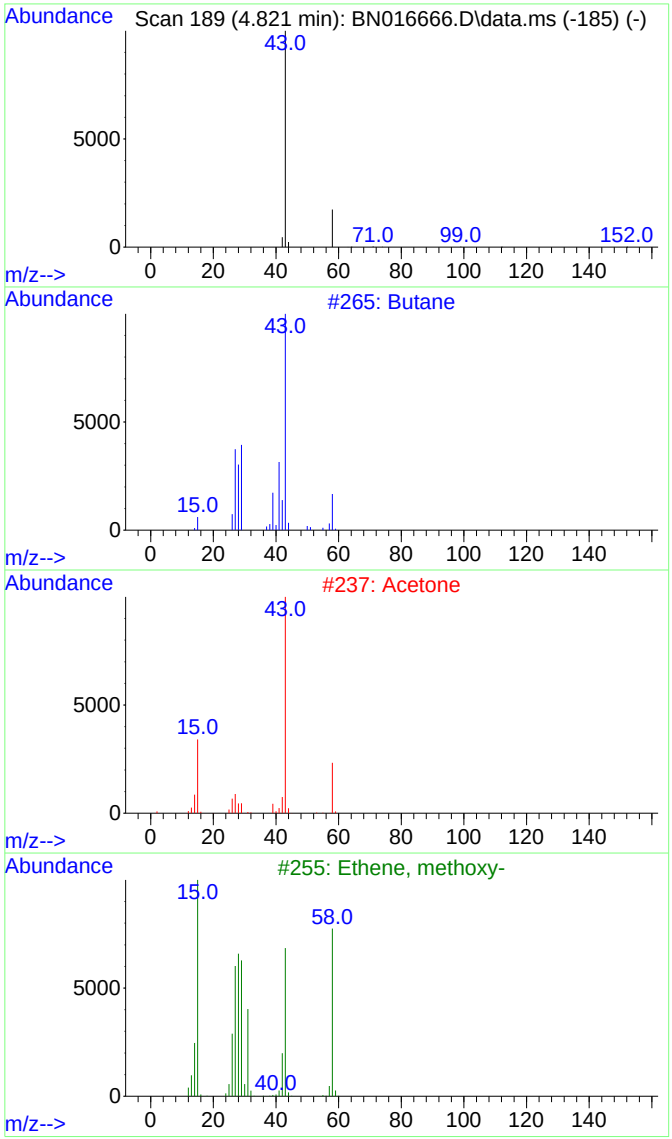
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

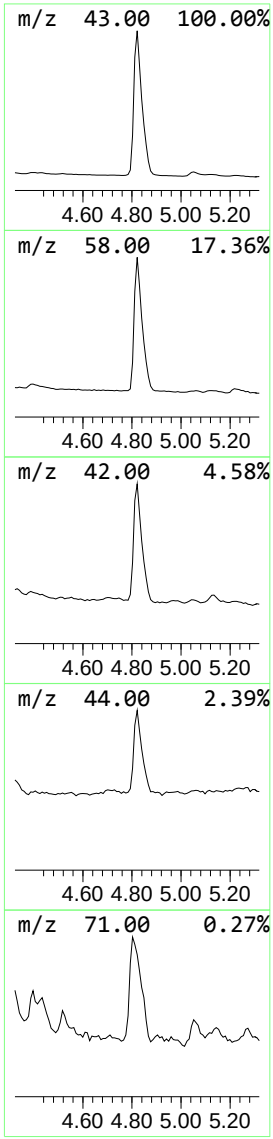
****

Peak Number 5 Butane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.821	0.43 ng	118327	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane	58	C4H10	000106-97-8	4
2	Acetone	58	C3H6O	000067-64-1	4
3	Ethene, methoxy-	58	C3H6O	000107-25-5	3
4	Diazene, dimethyl-	58	C2H6N2	000503-28-6	3
5	Hydrogen isocyanate	43	CHNO	000075-13-8	2





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ALS Vial : 34 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

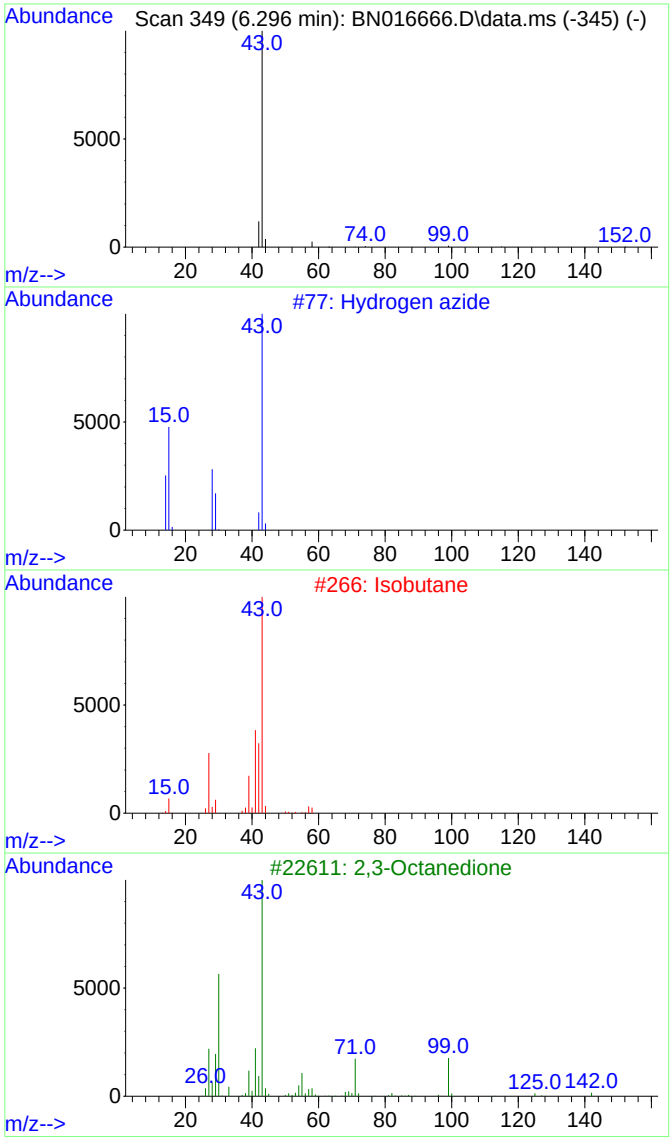
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

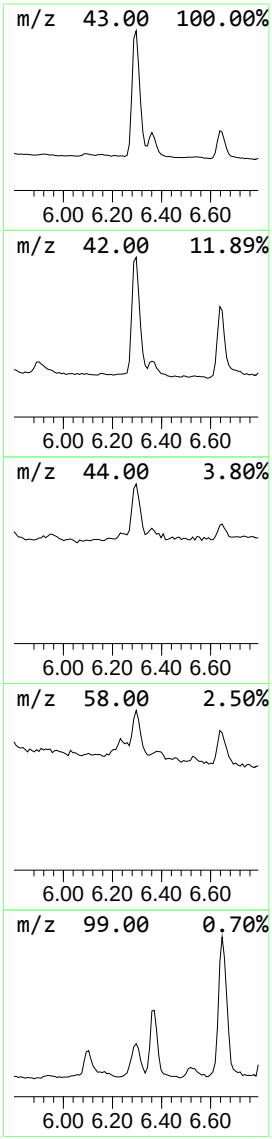
****

Peak Number 6 Hydrogen azide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.296	0.12 ng	31815	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrogen azide	43	HN3	007782-79-8	4
2	Isobutane	58	C4H10	000075-28-5	3
3	2,3-Octanedione	142	C8H14O2	000585-25-1	2
4	Hydrogen isocyanate	43	CHNO	000075-13-8	2
5	Carbonic acid, dipropyl ester	146	C7H14O3	000623-96-1	2





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Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

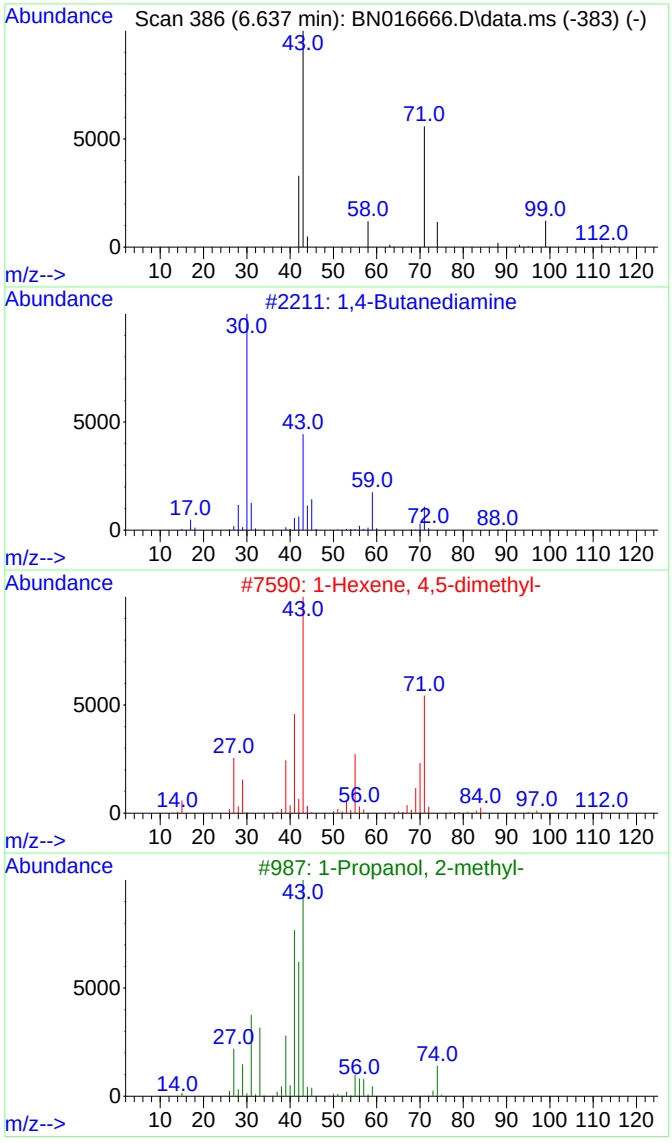
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

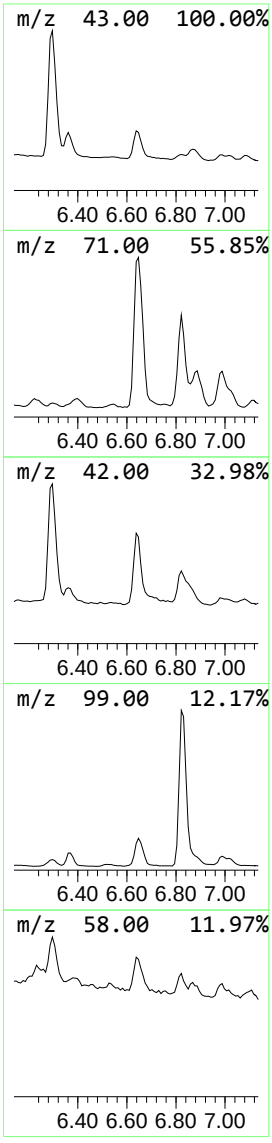
****

Peak Number 7 1,4-Butanediamine Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.637	0.06 ng	15832	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Butanediamine	88	C4H12N2	000110-60-1	5
2	1-Hexene, 4,5-dimethyl-	112	C8H16	016106-59-5	5
3	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	5
4	Acetic acid, hydrazide	74	C2H6N2O	001068-57-1	4
5	2-Propanone, 1-hydroxy-	74	C3H6O2	000116-09-6	4





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BNA_N
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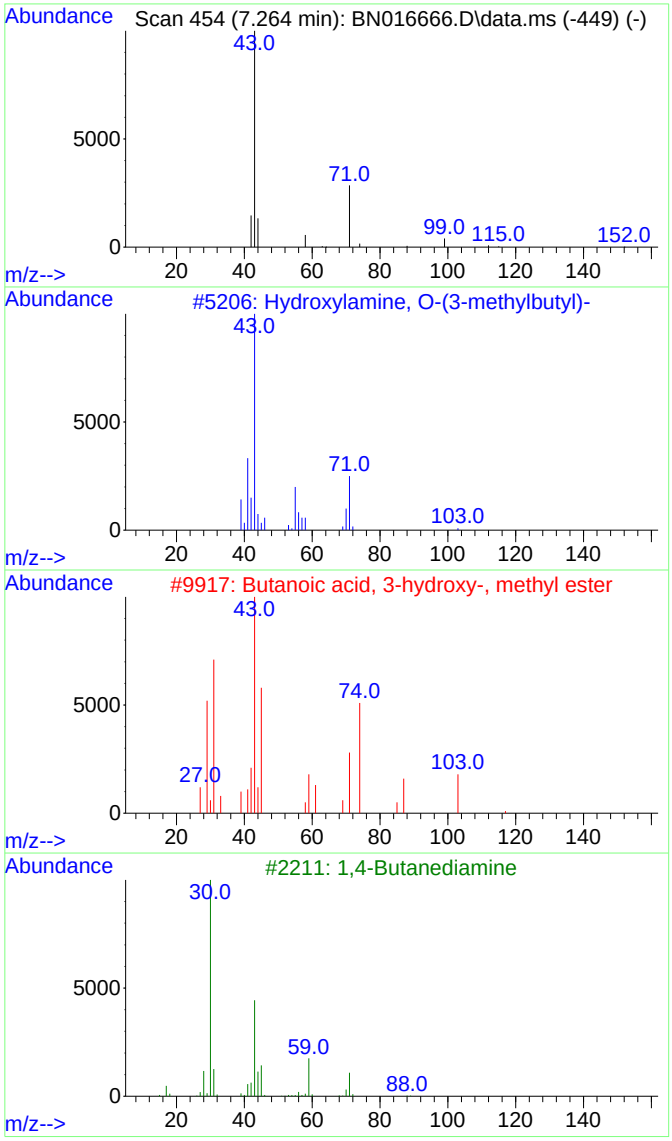
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

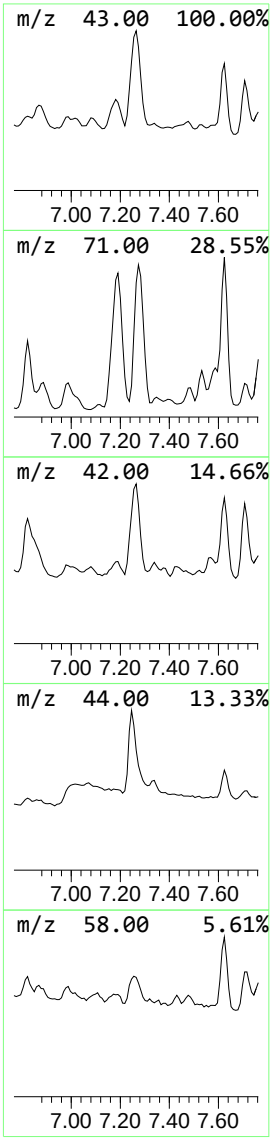
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Hydroxylamine, O-(3-methylb... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.264	0.08 ng	21739	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydroxylamine, O-(3-methylbutyl)-	103	C5H13NO	019411-65-5	9
2	Butanoic acid, 3-hydroxy-, methy...	118	C5H10O3	001487-49-6	9
3	1,4-Butanediamine	88	C4H12N2	000110-60-1	7
4	Hex-4-enylamine	99	C6H13N	055108-01-5	5
5	Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	4





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Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

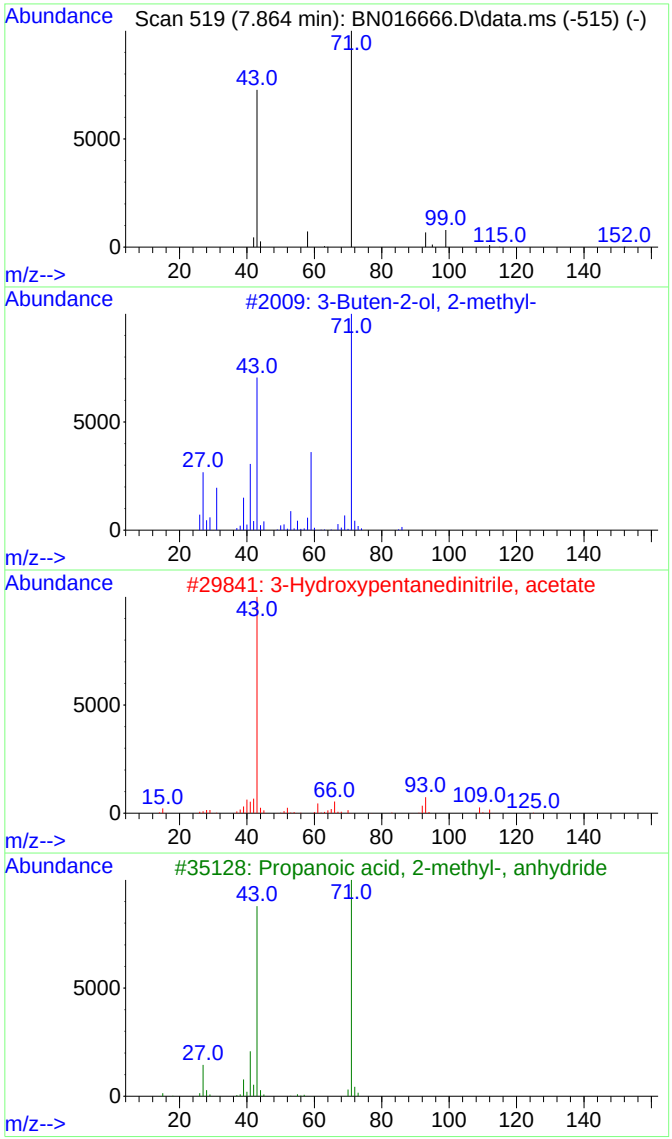
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

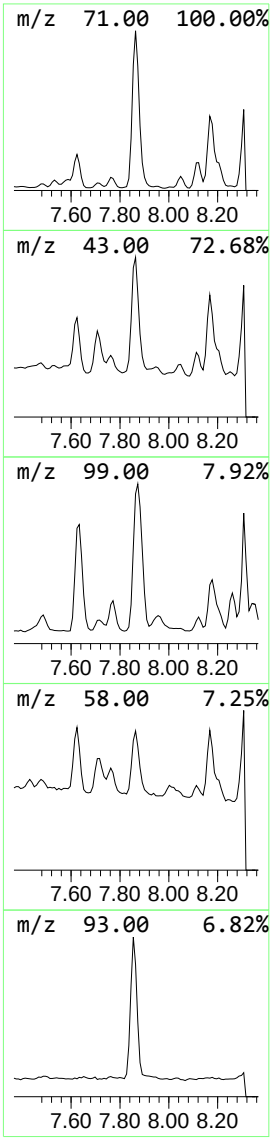
****

Peak Number 9 3-Buten-2-ol, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.864	0.14 ng	39159	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	4
2	3-Hydroxypentanedinitrile, acetate	152	C7H8N2O2	1000506-60-2	4
3	Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	4
4	Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	4
5	3-Penten-2-ol	86	C5H10O	001569-50-2	4





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Data File : BN016666.D
Acq On : 02 Oct 2021 11:08
Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : RE114D1-20210916

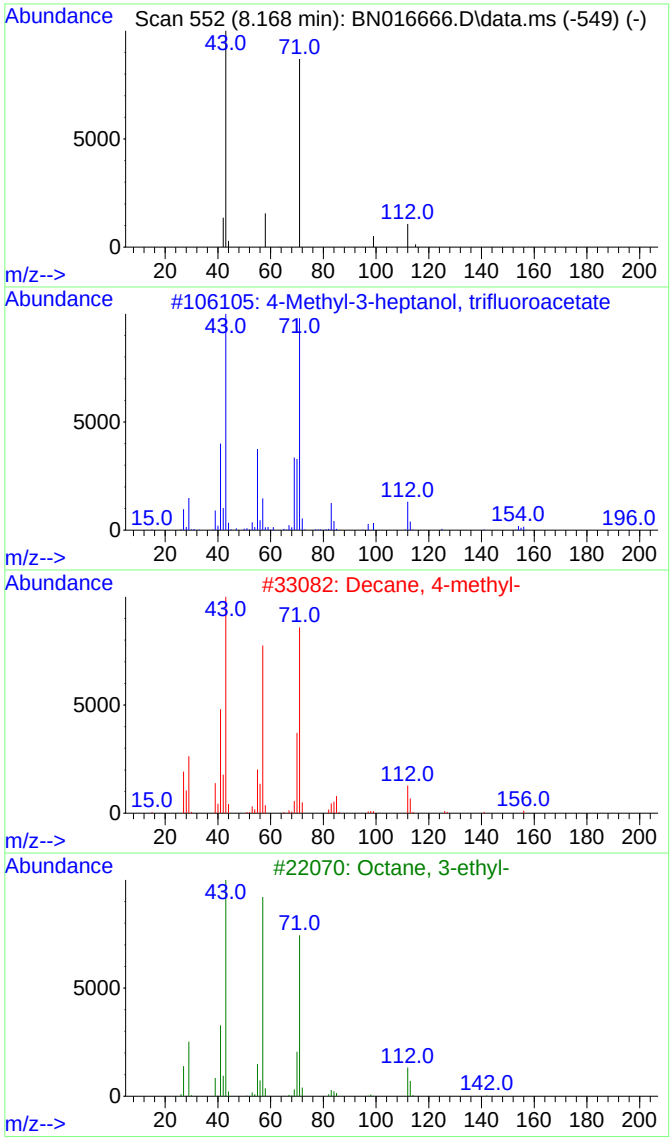
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

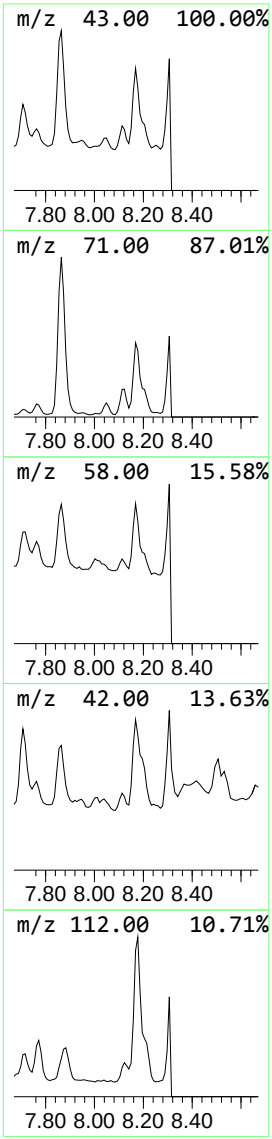
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 4-Methyl-3-heptanol, triflu... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.168	0.11 ng	29002	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	43
2	Decane, 4-methyl-	156	C11H24	002847-72-5	9
3	Octane, 3-ethyl-	142	C10H22	005881-17-4	9
4	1,3-Butadien-1-ol, acetate	112	C6H8O2	001515-76-0	7
5	Butane	58	C4H10	000106-97-8	7





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Operator : CG/JU
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Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

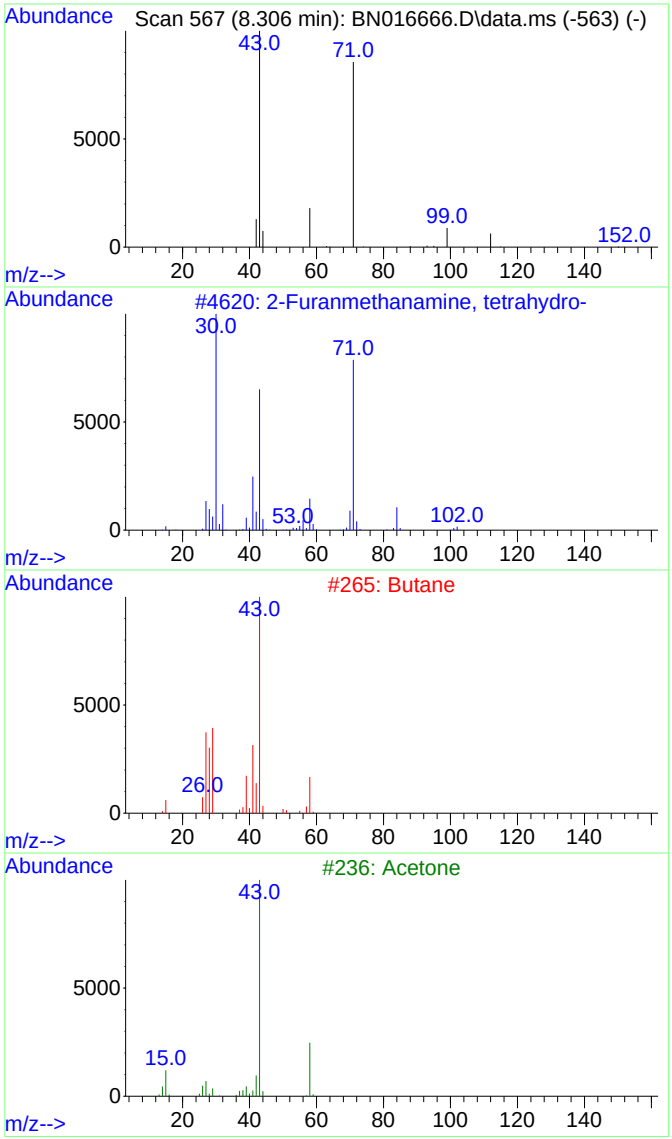
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

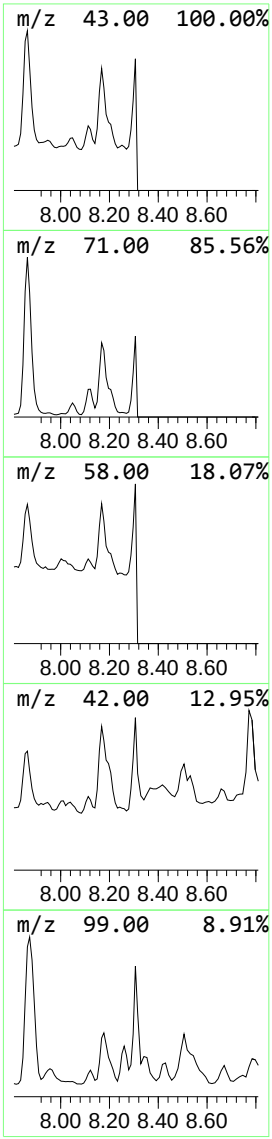
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 2-Furanmethanamine, tetrahy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.306	0.08 ng	20753	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanmethanamine, tetrahydro-	101	C5H11NO	004795-29-3	9
2	Butane	58	C4H10	000106-97-8	5
3	Acetone	58	C3H6O	000067-64-1	5
4	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	4
5	1,4-Butanediamine	88	C4H12N2	000110-60-1	4





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Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

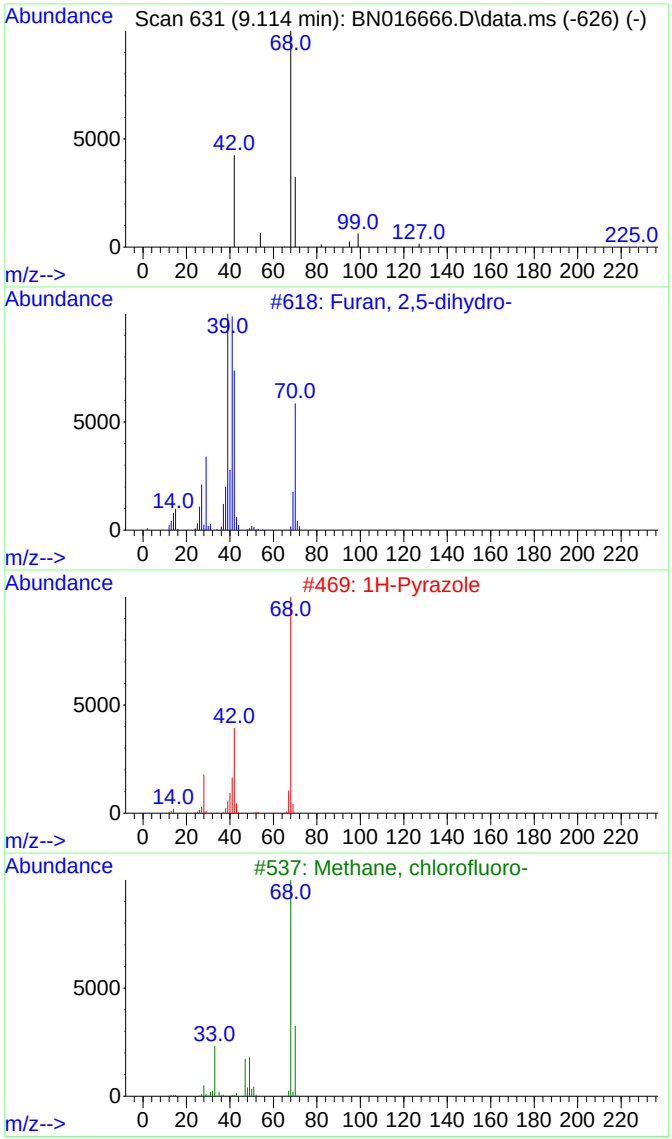
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

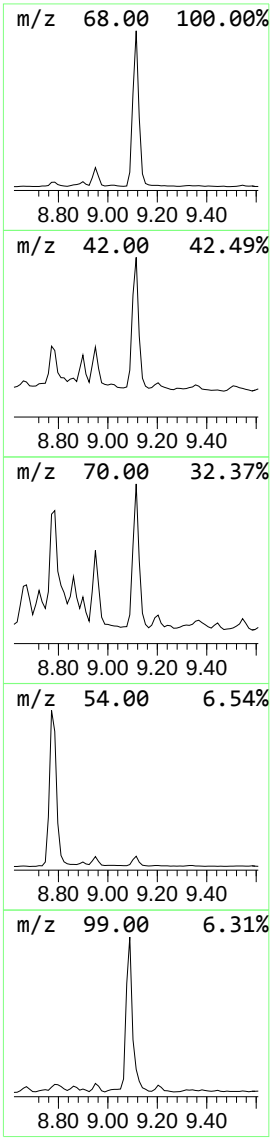
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Furan, 2,5-dihydro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.114	0.06 ng	25677	Naphthalene-d8	10.407

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	4
2	1H-Pyrazole	68	C3H4N2	000288-13-1	4
3	Methane, chlorofluoro-	68	CH2ClF	000593-70-4	4
4	1,4-Pentadiene	68	C5H8	000591-93-5	4
5	Acetic acid, cyano-, methyl ester	99	C4H5NO2	000105-34-0	4





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Operator : CG/JU
Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

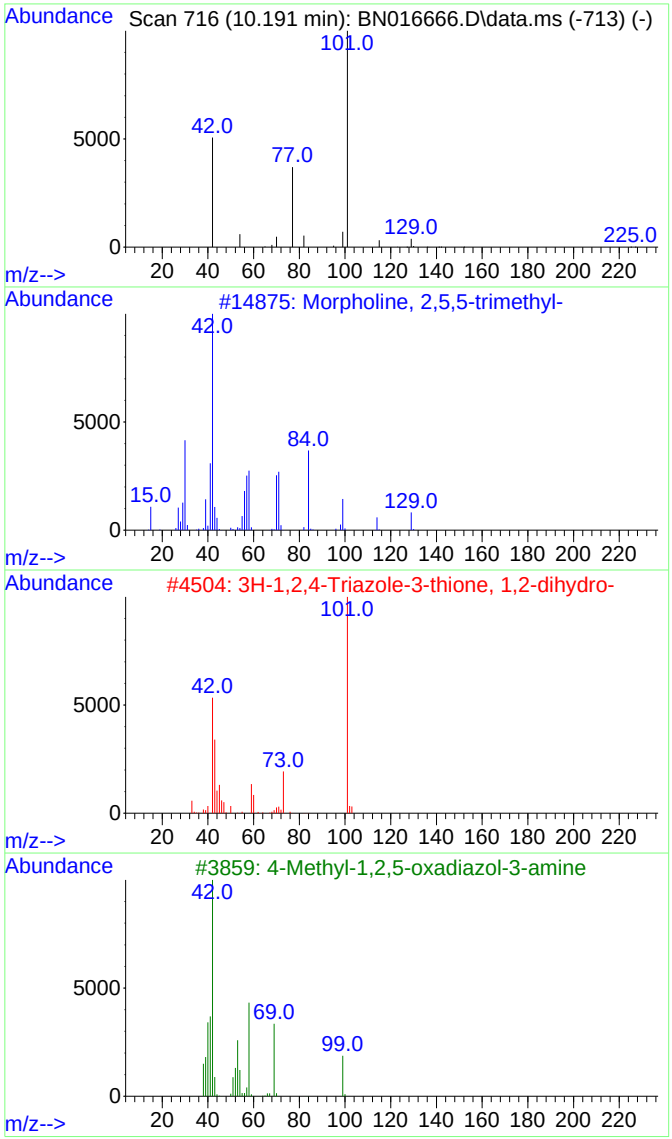
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

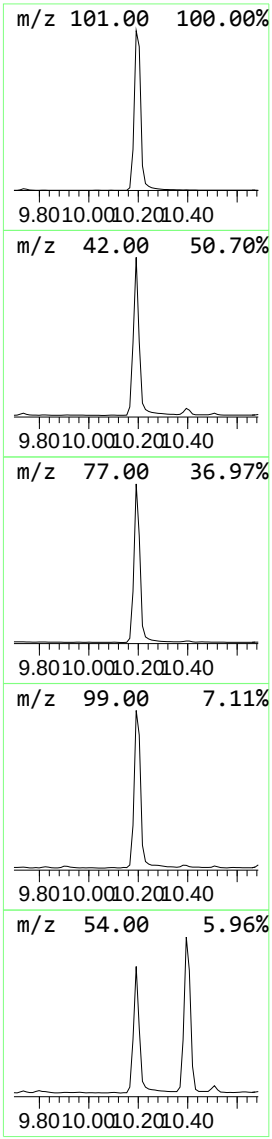
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Morpholine, 2,5,5-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.191	0.70 ng	319338	Naphthalene-d8	10.407

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Morpholine, 2,5,5-trimethyl-	129	C7H15NO	1000460-73-8	5
2	3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	5
3	4-Methyl-1,2,5-oxadiazol-3-amine	99	C3H5N3O	1000411-31-0	5
4	Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	5
5	Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	5





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Sample : M3829-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

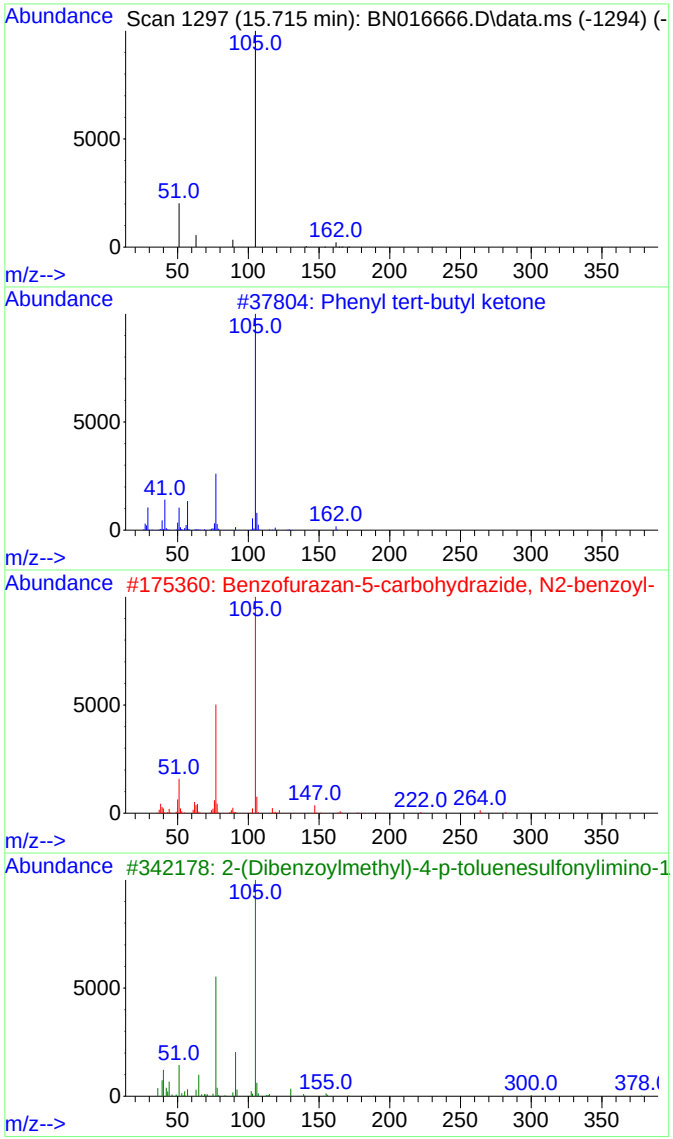
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

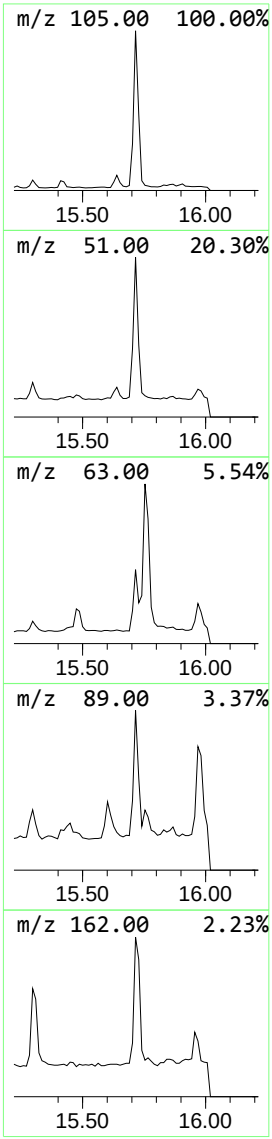
****

Peak Number 14 Phenyl tert-butyl ketone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.715	0.05 ng	21438	Phenanthrene-d10	17.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenyl tert-butyl ketone	162	C11H14O	000938-16-9	3
2	Benzofurazan-5-carbohydrazide, N...	282	C14H10N4O3	1000272-94-4	2
3	2-(Dibenzoylmethyl)-4-p-toluenes...	533	C32H23NO5S	300360-29-6	2
4	.alpha.,.alpha.-Dichloroacetophe...	188	C8H6Cl2O	002648-61-5	2
5	Propanoic acid, 2-(benzoylamino)...	283	C17H17NO3	074923-17-4	2





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Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D1-20210916

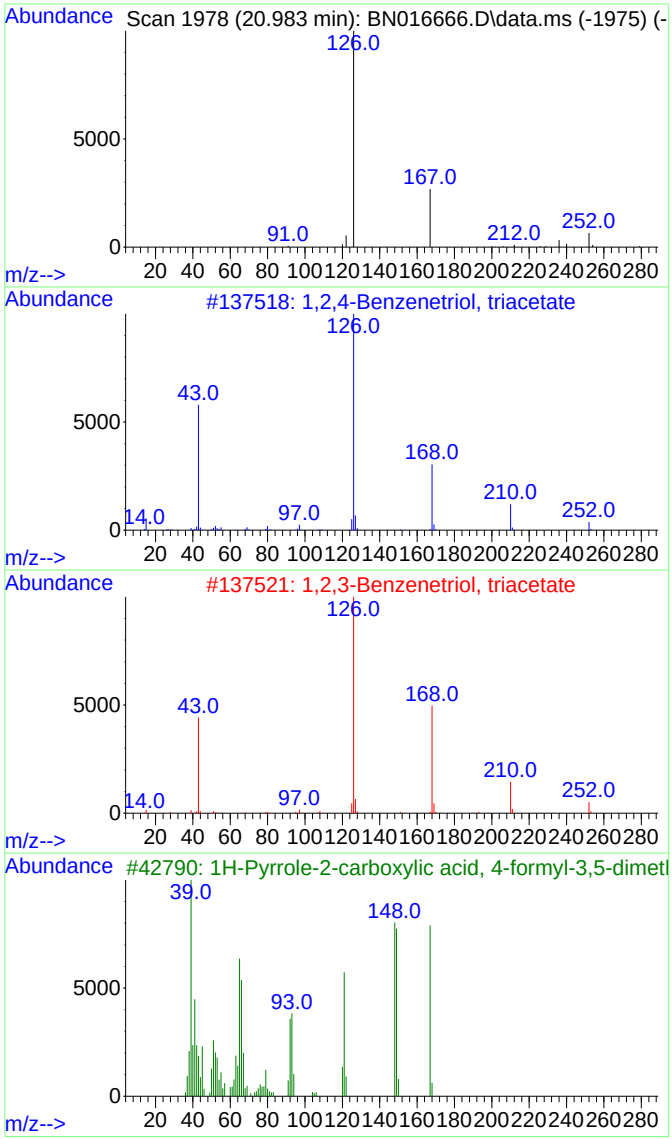
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

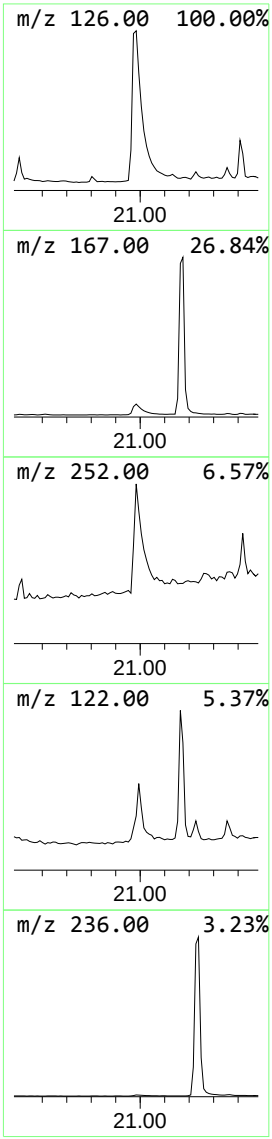
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1,2,4-Benzenetriol, triacetate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.983	0.04 ng	20711	Chrysene-d12	21.238

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4-Benzenetriol, triacetate	252	C12H12O6	000613-03-6	4
2	1,2,3-Benzenetriol, triacetate	252	C12H12O6	000525-52-0	4
3	1H-Pyrrole-2-carboxylic acid, 4-...	167	C8H9NO3	1000362-61-5	4
4	O-Ethyl N,N-dimethyl-P-methylpho...	167	C5H14NOPS	276694-08-7	4
5	4-Morpholinopiperidine, N-(2-met...	240	C13H24N2O2	1010469-51-1	4





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Sample : M3829-02
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Instrument :
BNA_N
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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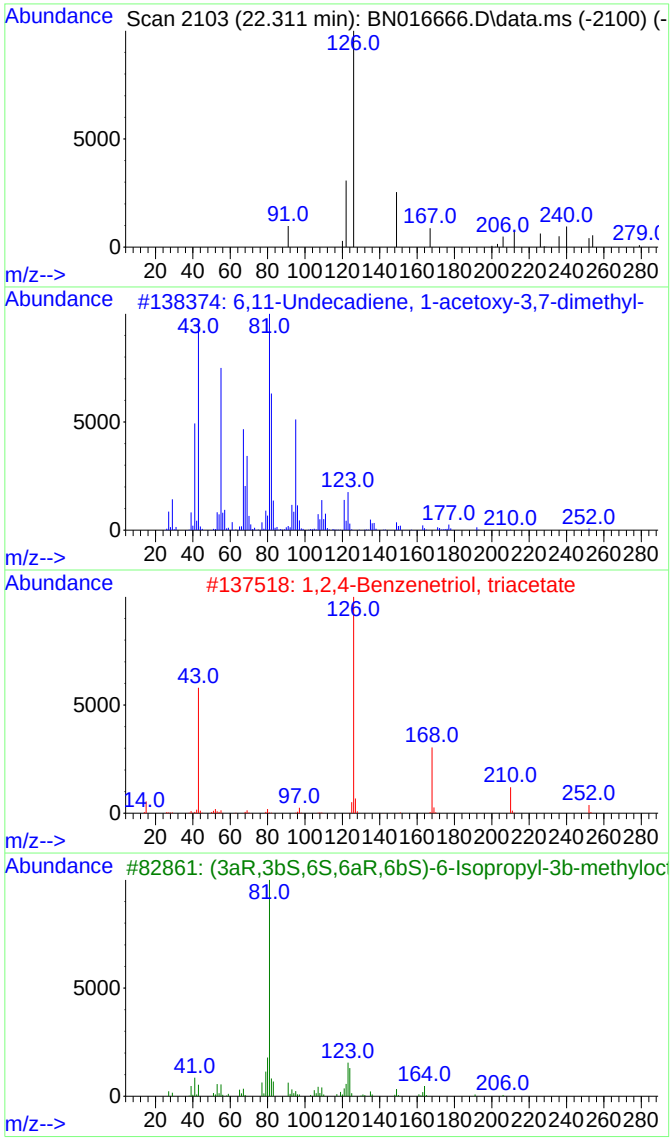
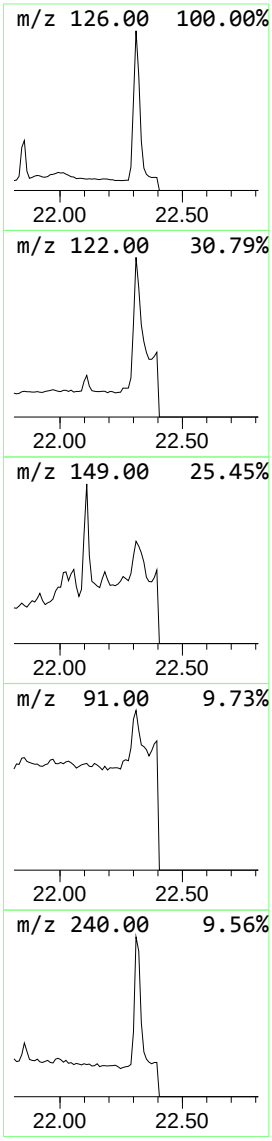
****
Peak Number 16 6,11-Undecadiene, 1-acetoxy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.311	0.04 ng	22226	Chrysene-d12	21.238

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6,11-Undecadiene, 1-acetoxy-3,7-...	252	C16H28O2	1000150-66-0	4
2	1,2,4-Benzenetriol, triacetate	252	C12H12O6	000613-03-6	4
3	(3aR,3bS,6S,6aR,6bS)-6-Isopropyl...	206	C14H22O	013844-03-6	4
4	1,2,3-Benzenetriol, triacetate	252	C12H12O6	000525-52-0	4
5	(6S,12aR,E)-2,2-Dimethyltetrahyd...	236	C14H24N2O	155975-43-2	4





Tentatively Identified Compound (LSC) summary

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					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Pyridine, 1-oxide	3.161	4.1	ng	1128600	1	7.643	108847	0.4
Hydrogen isocya...	3.364	0.9	ng	232441	1	7.643	108847	0.4
Acetic acid, me...	3.816	0.1	ng	34586	1	7.643	108847	0.4
Acetamide, N-et...	4.010	0.1	ng	25869	1	7.643	108847	0.4
Butane	4.821	0.4	ng	118327	1	7.643	108847	0.4
Hydrogen azide	6.296	0.1	ng	31815	1	7.643	108847	0.4
1,4-Butanediamine	6.637	0.1	ng	15832	1	7.643	108847	0.4
Hydroxylamine, ...	7.264	0.1	ng	21739	1	7.643	108847	0.4
3-Buten-2-ol, 2...	7.864	0.1	ng	39159	1	7.643	108847	0.4
4-Methyl-3-hept...	8.168	0.1	ng	29002	1	7.643	108847	0.4
2-Furanmethanam...	8.306	0.1	ng	20753	1	7.643	108847	0.4
Furan, 2,5-dihy...	9.114	0.1	ng	25677	2	10.407	182621	0.4
Morpholine, 2,5...	10.191	0.7	ng	319338	2	10.407	182621	0.4
Phenyl tert-but...	15.715	0.1	ng	21438	4	17.018	167746	0.4
1,2,4-Benzenetr...	20.983	0.0	ng	20711	5	21.238	202138	0.4
6,11-Undecadien...	22.311	0.0	ng	22226	5	21.238	202138	0.4