LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : MW202D1-20210919

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: BN016642.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.502	44	46	59	rBV	12166	42930	13.62%	1.474%
2	3.825	77	81	98	rVB2	13273	48271	15.31%	1.657%
3	4.821	186	189	196	rVB	67202	134796	42.76%	4.628%
4	5.052	211	214	220	rBV	3013	7192	2.28%	0.247%
5	5.273	234	238	244	rVV	9373	20585	6.53%	0.707%
6	5.365	245	248	257	rVB	2995	7886	2.50%	0.271%
7	6.656	383	388	391	rBV	9497	19353	6.14%	0.664%
8	6.850	403	409	420	rVB2	11382	34978	11.10%	1.201%
9	7.265	449	454	460	rVB3	9074	25636	8.13%	0.880%
10	7.643	490	495	500	rBV	56445	106635	33.83%	3.661%
11	7.707	500	502	506	rVB	4118	5845	1.85%	0.201%
12	7.864	515	519	525	rVB	9769	19790	6.28%	0.679%
13	8.177	549	553	555	rBV	5489	11213	3.56%	0.385%
14	8.306	565	567	569	rVB2	9619	8834	2.80%	0.303%
15	8.455	574	579	581	rBV2	3040	8821	2.80%	0.303%
16	8.785	602	605	610	rBV	33575	61780	19.60%	2.121%
17	8.949	616	618	621	rVB	4891	7767	2.46%	0.267%
18	9.748	678	681	685	rBV	4465	8034	2.55%	0.276%
19	10.204	713	717	720	rBV	182536	315213	100.00%	10.823%
20	10.407	730	733	736	rBV	106545	182736	57.97%	6.274%
21	10.774	759	762	768	rBV2	2927	8953	2.84%	0.307%
22	11.205	793	796	802	rVB2	4885	9699	3.08%	0.333%
23	11.307	802	804	814	rVB	3259	7812	2.48%	0.268%
24	12.007	915	932	942	rBV	51993	84157	26.70%	2.890%
25	12.899	1071	1075	1078	rVB2	101225	188838	59.91%	6.484%
26	13.114	1089	1092	1095	rBV	6154	10034	3.18%	0.345%
27	13.178	1095	1097	1101	rVV2	12037	23511	7.46%	0.807%
28	13.812	1144	1147	1153	rVB3	4381	9687	3.07%	0.333%
29	14.269	1180	1183	1186	rBV	138973	202400	64.21%	6.949%
30	14.434	1193	1196	1199	rBV	11917	17595	5.58%	0.604%
31	14.523	1199	1203	1207	rVB3	2995	7432	2.36%	0.255%
32	15.639	1289	1291	1295	rVB2	4371	8217	2.61%	0.282%
33	15.715	1295	1297	1299	rBV	13674	21845	6.93%	0.750%
34	15.766	1299	1301	1304	rVB2	17389	27205	8.63%	0.934%
35	15.969	1311	1317	1321	rVB3	5501	26888	8.53%	0.923%
36	16.835	1383	1388	1391	rBV3	13674	25867	8.21%	0.888%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : MW202D1-20210919

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

37	17.018	1400	1403	1406	rBV	97114	167423	53.11%	5.749%
38	17.724	1458	1461	1465	rBV2	3171	7836	2.49%	0.269%
39	17.955	1477	1480	1483	rBV	14457	24932	7.91%	0.856%
40	18.909	1661	1665	1672	rVB	4824	5522	1.75%	0.190%
41	19.063	1690	1696	1718	rBV	136981	189287	60.05%	6.499%
42	19.247	1728	1733	1746	rBV	15176	22049	6.99%	0.757%
43	19.679	1814	1820	1832	rBV	131627	161344	51.19%	5.540%
44	20.017	1885	1887	1889	rBV2	13078	28484	9.04%	0.978%
45	20.378	1919	1921	1923	rBV	5515	6857	2.18%	0.235%
46	20.973	1975	1977	1980	rBV	11666	17979	5.70%	0.617%
47	21.026	1980	1982	1987	rVB	13022	17004	5.39%	0.584%
48	21.164	1993	1995	1998	rBV	46882	65060	20.64%	2.234%
49	21.228	1998	2001	2007	rVB	136621	199965	63.44%	6.866%
50	22.109	2081	2084	2089	rVB	11550	16774	5.32%	0.576%
51	23.420	2403	2408	2414	rVB	4630	5895	1.87%	0.202%
52	23.485	2415	2427	2451	rVB	98043	187675	59.54%	6.444%
53	24.084	2595	2602	2617	rVB	6306	11488	3.64%	0.394%
54	24.854	2820	2827	2842	rVB	5594	11866	3.76%	0.407%
55	25.757	3080	3091	3112	rVB3	2283	6566	2.08%	0.225%

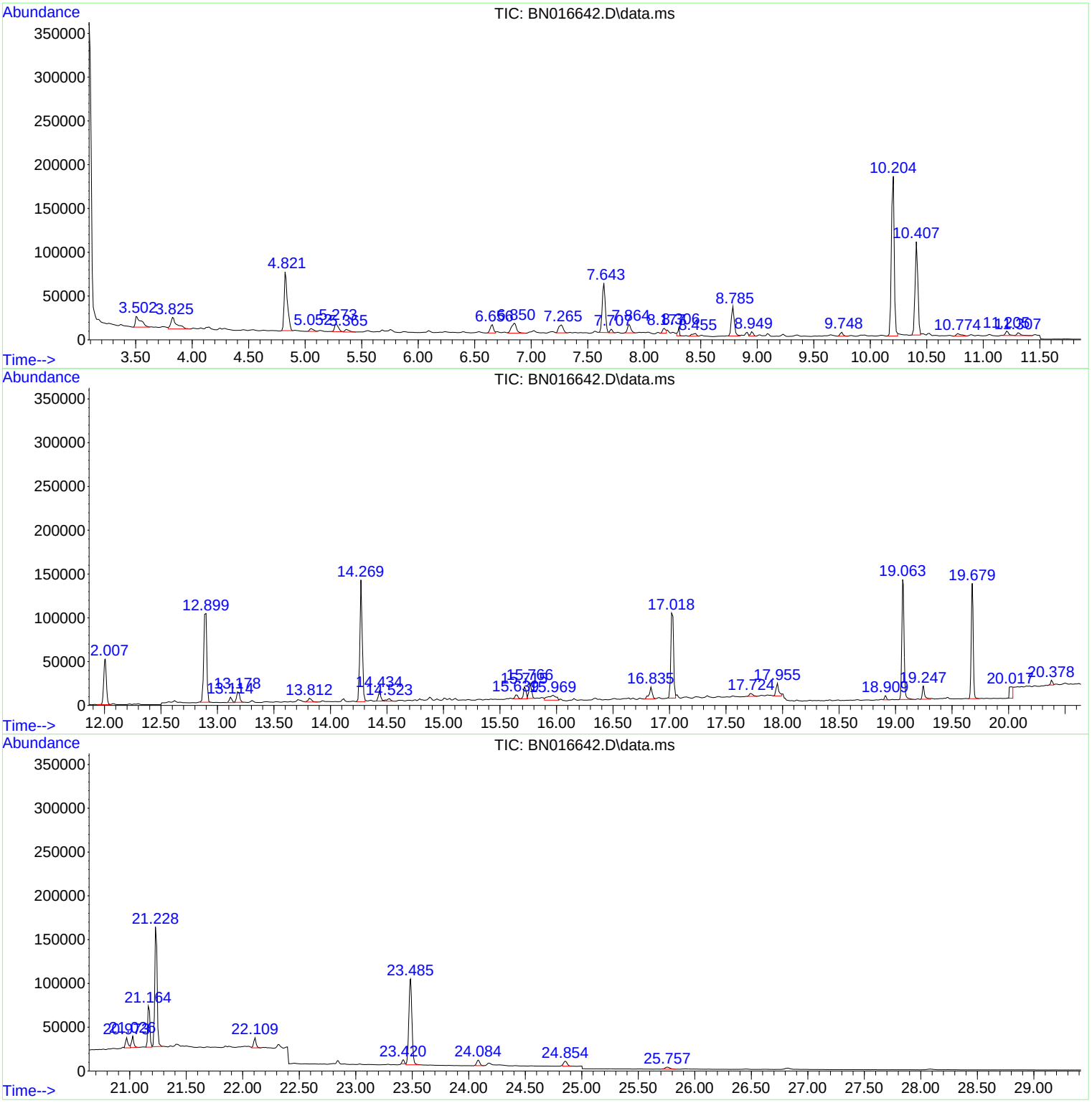
Sum of corrected areas: 2912441

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

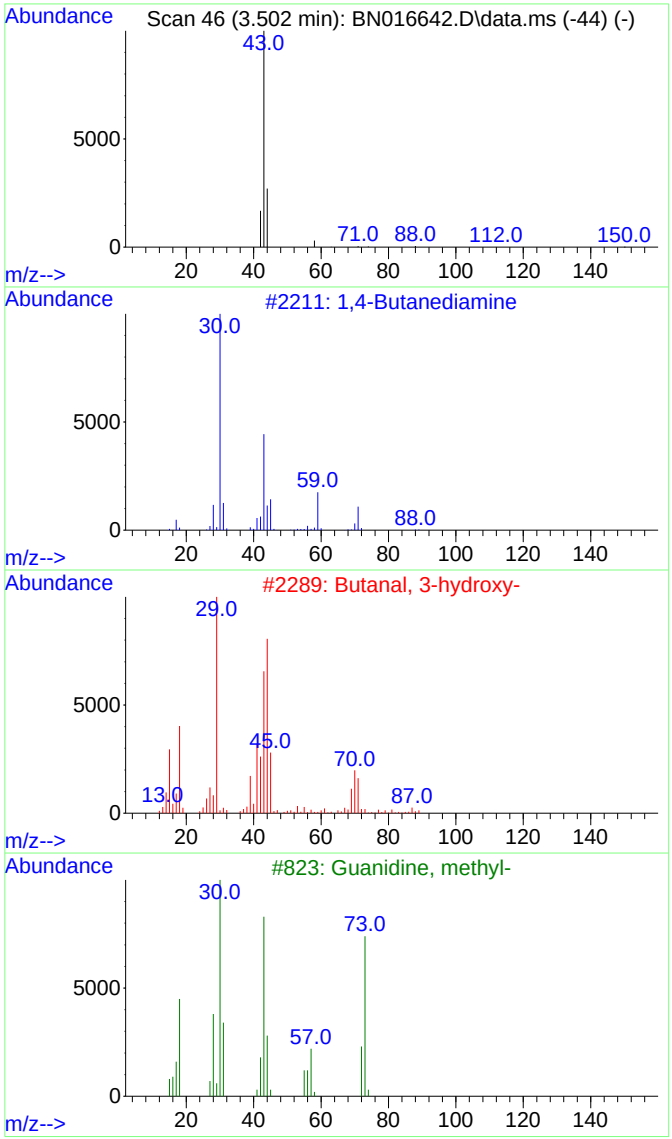
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

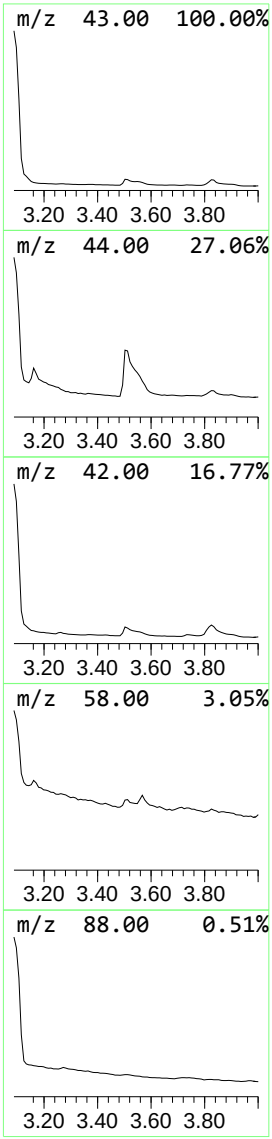
****

Peak Number 1 1,4-Butanediamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.502	0.16 ng	42930	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	Butanal, 3-hydroxy-	88	C4H8O2	000107-89-1	4
3	Guanidine, methyl-	73	C2H7N3	000471-29-4	4
4	2-Amino-2-oxoethyl acetate	117	C4H7NO3	1000500-76-5	4
5	Hydrogen isocyanate	43	CHNO	000075-13-8	3





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : MW202D1-20210919

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

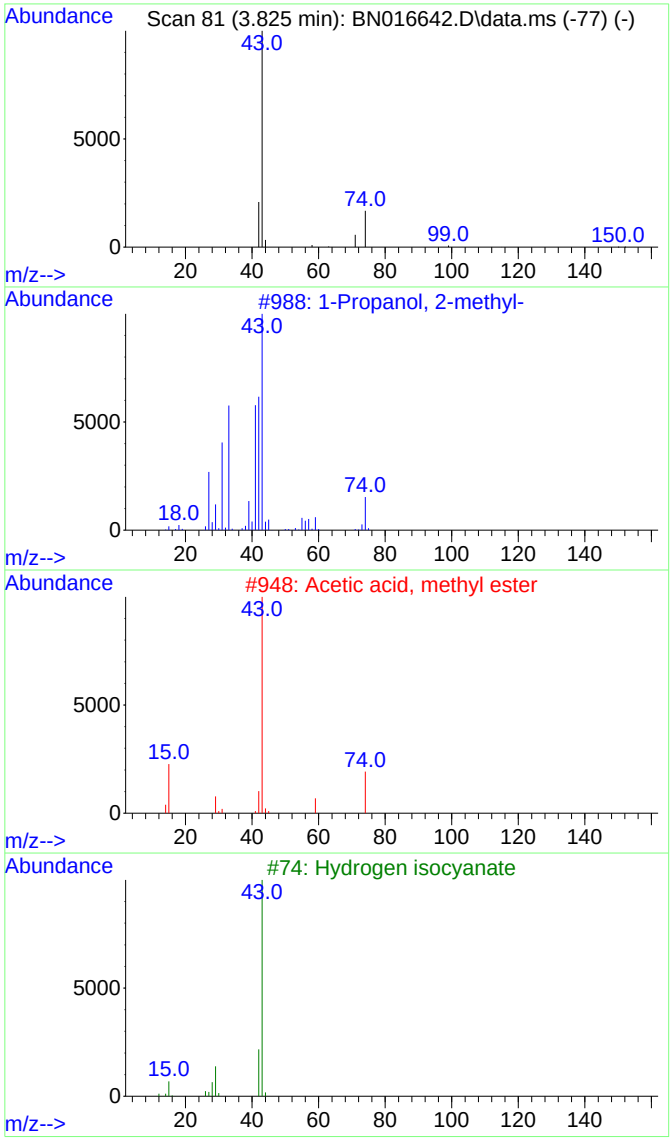
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

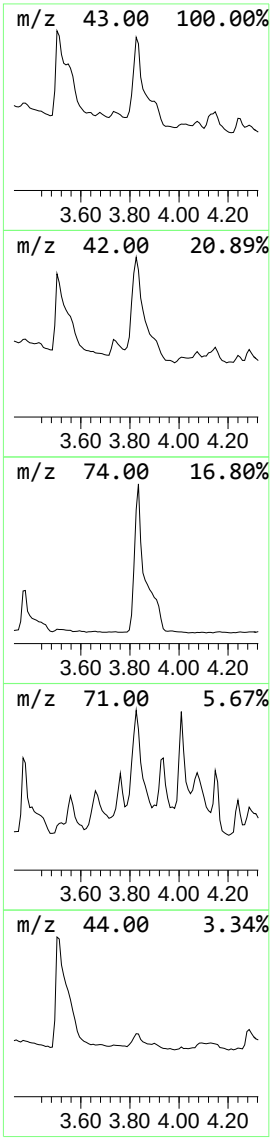
****

Peak Number 2 1-Propanol, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.825	0.18 ng	48271	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	4
2	Acetic acid, methyl ester	74	C3H6O2	000079-20-9	4
3	Hydrogen isocyanate	43	CHNO	000075-13-8	4
4	2-Propanone, 1-hydroxy-	74	C3H6O2	000116-09-6	4
5	Acetic acid, hydrazide	74	C2H6N2O	001068-57-1	3





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

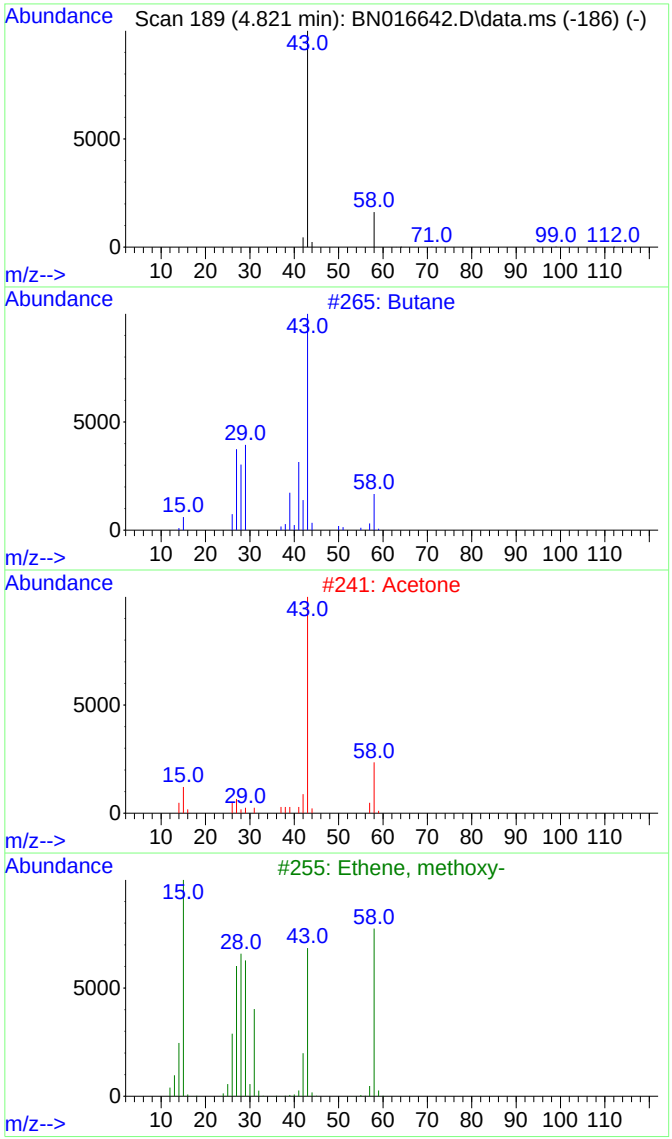
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

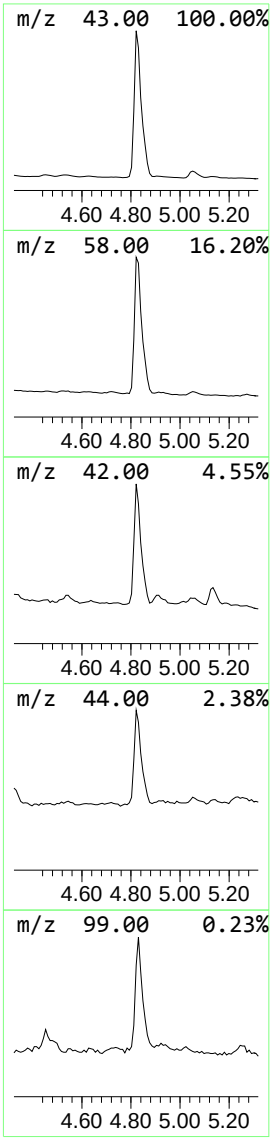
****

Peak Number 3 Butane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.821	0.51 ng	134796	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





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

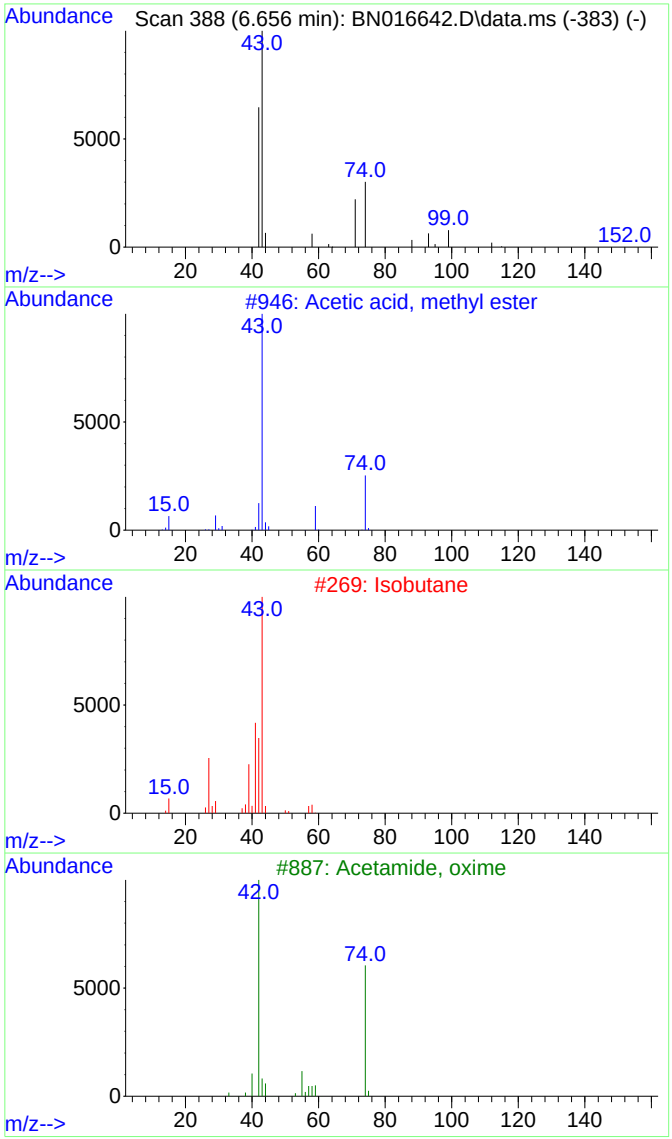
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

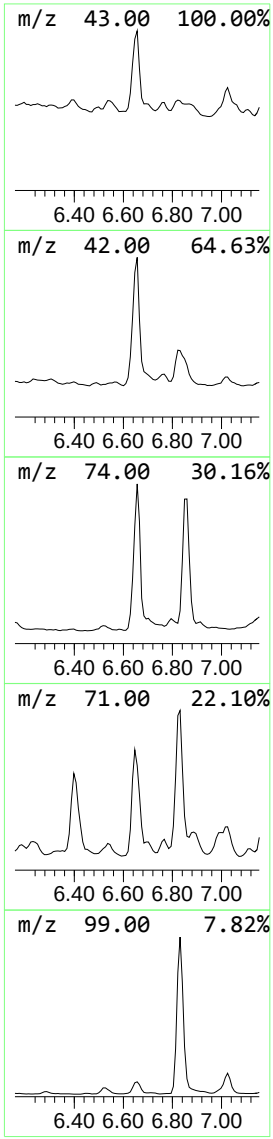
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Acetic acid, methyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.656	0.07 ng	19353	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	5
2	Isobutane	58	C4H10	000075-28-5	4
3	Acetamide, oxime	74	C2H6N2O	022059-22-9	4
4	Hydroxylamine, O-(2-methylpropyl)-	89	C4H11NO	005618-62-2	4
5	Pentane, 1-(2-propenyloxy)-	128	C8H16O	023186-70-1	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

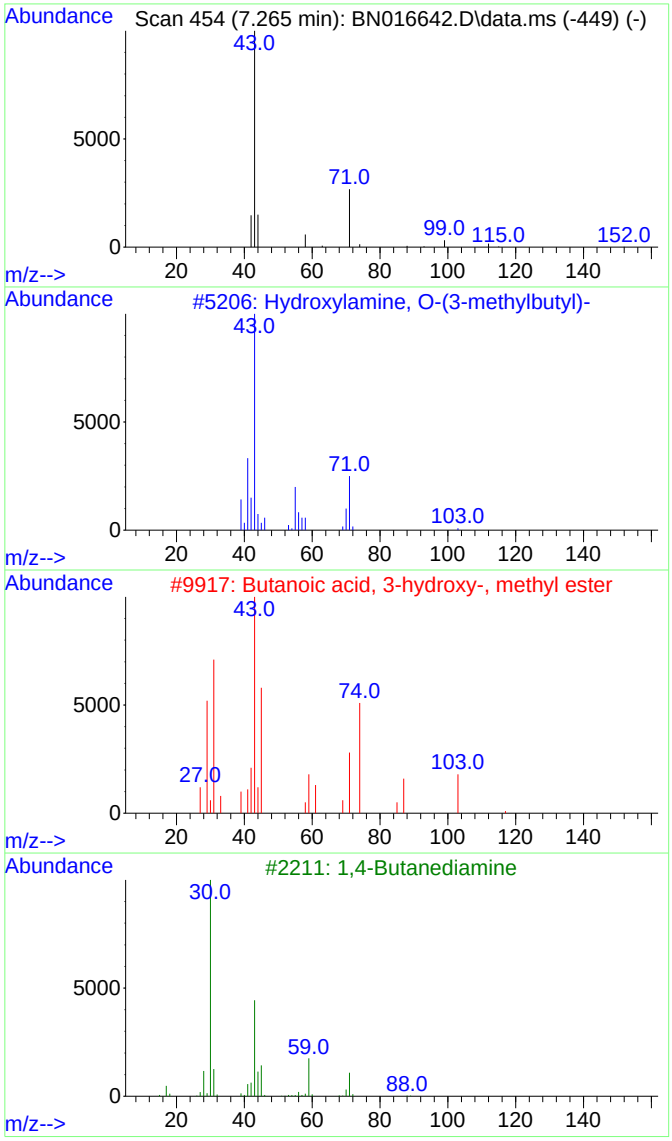
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

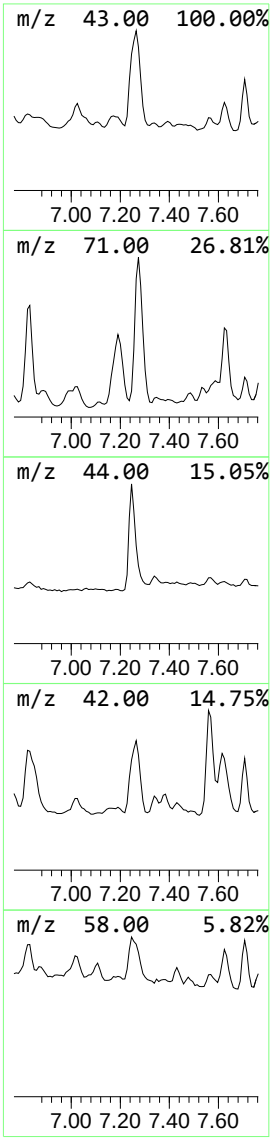
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hydroxylamine, O-(3-methylb... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.265	0.10 ng	25636	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	N-(2-Aminoethyl)-N-methylmethane...	152	C4H12N2O2S	778572-84-2	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

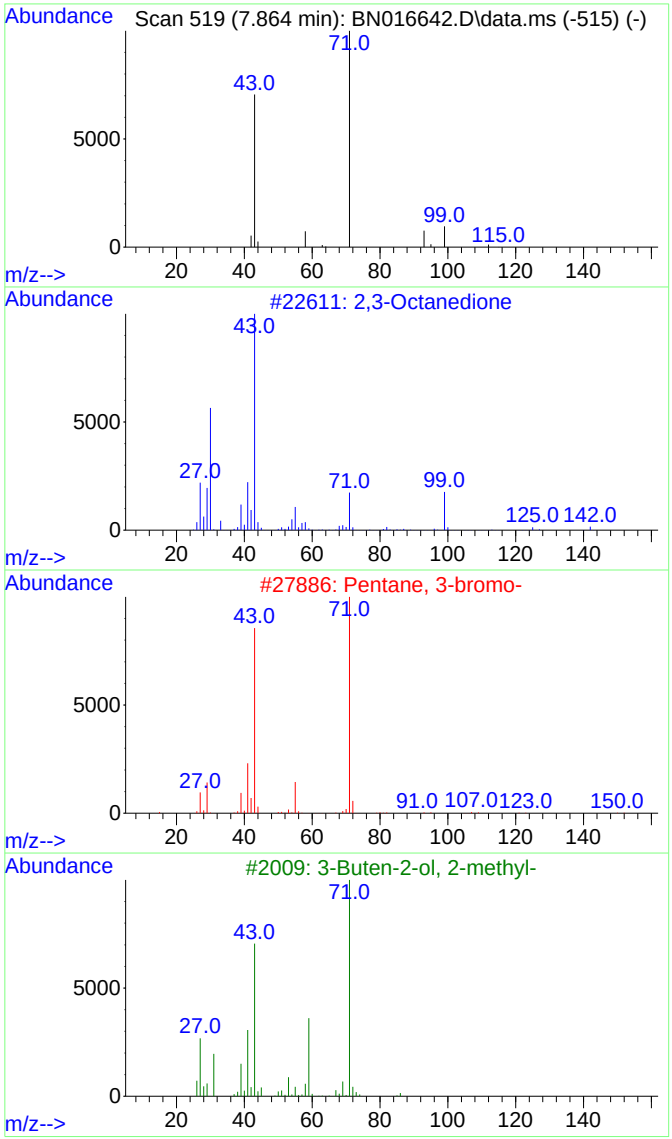
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

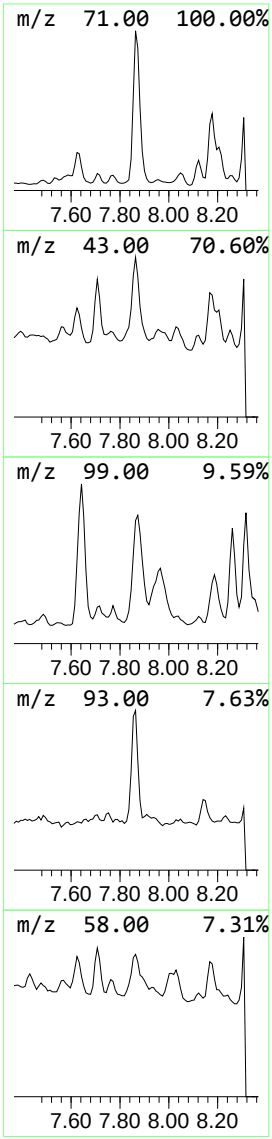
****

Peak Number 6 2,3-Octanedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.864	0.07 ng	19790	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Octanedione	142	C8H14O2	000585-25-1	4
2	Pentane, 3-bromo-	150	C5H11Br	001809-10-5	4
3	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	4
4	Methallyl butyrate	142	C8H14O2	1000428-83-8	4
5	Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	4





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Data File : BN016642.D
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Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

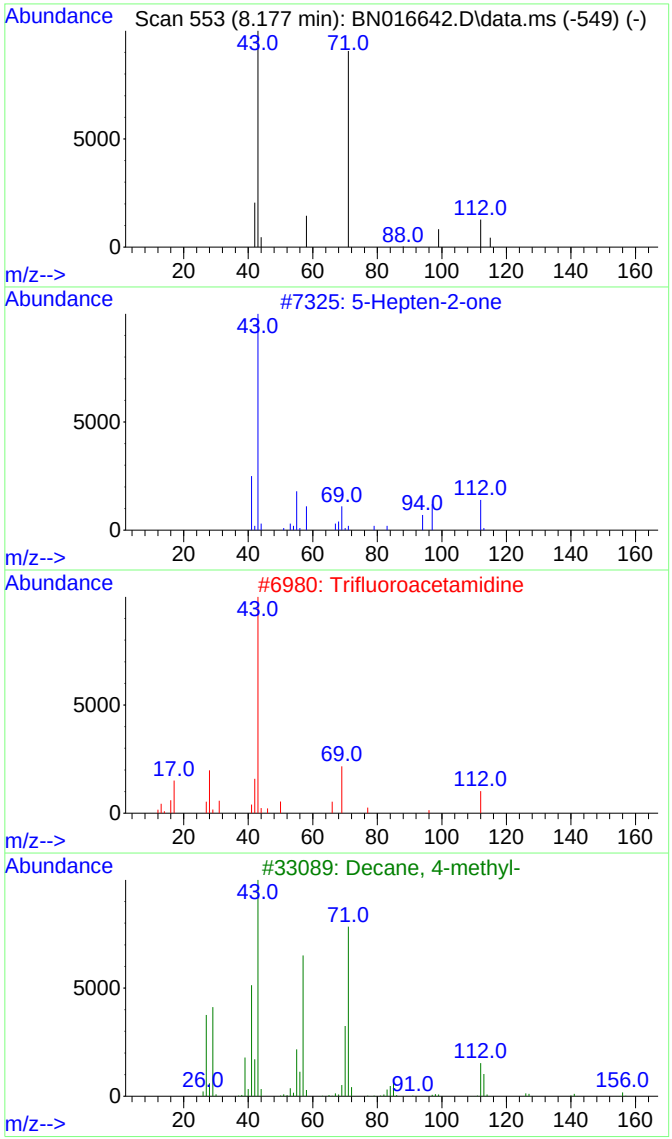
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

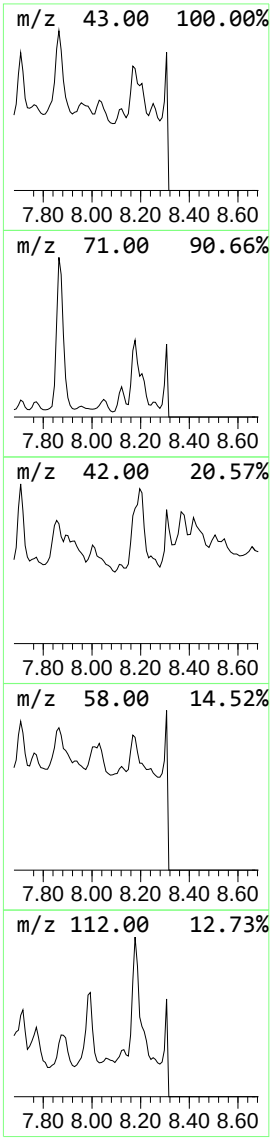
****

Peak Number 7 5-Hepten-2-one Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.177	0.04 ng	11213	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Hepten-2-one	112	C7H12O	006714-00-7	9
2	Trifluoroacetamidine	112	C2H3F3N2	000354-37-0	9
3	Decane, 4-methyl-	156	C11H24	002847-72-5	9
4	1-Trifluoroacetoxy-2-methylpentane	198	C8H13F3O2	155089-96-6	9
5	n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

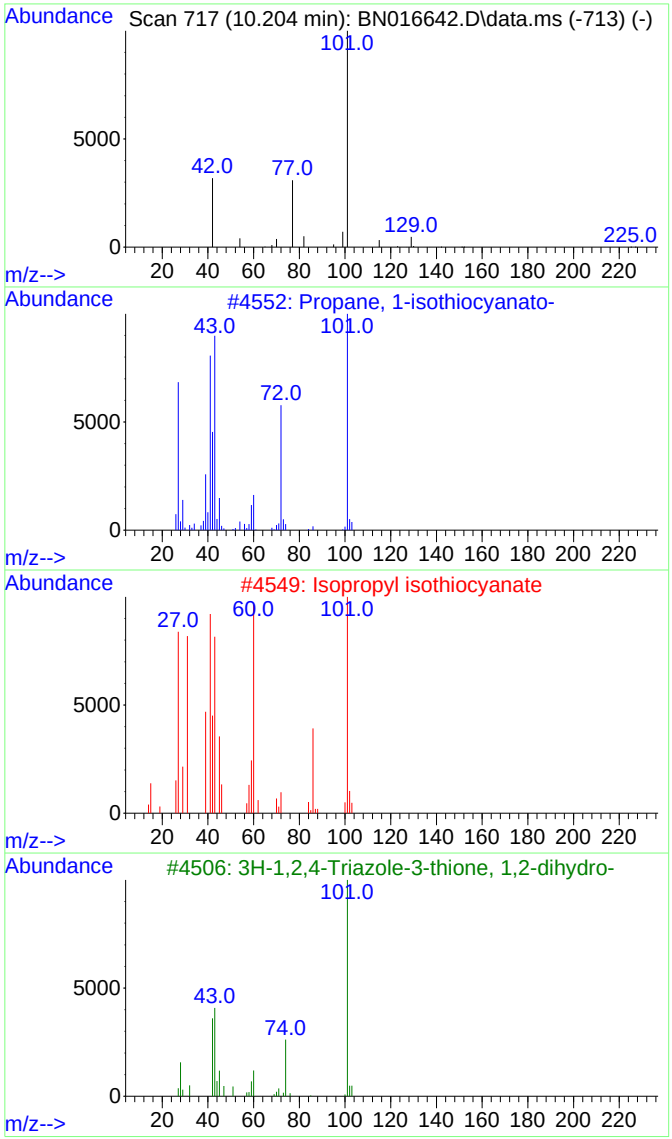
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

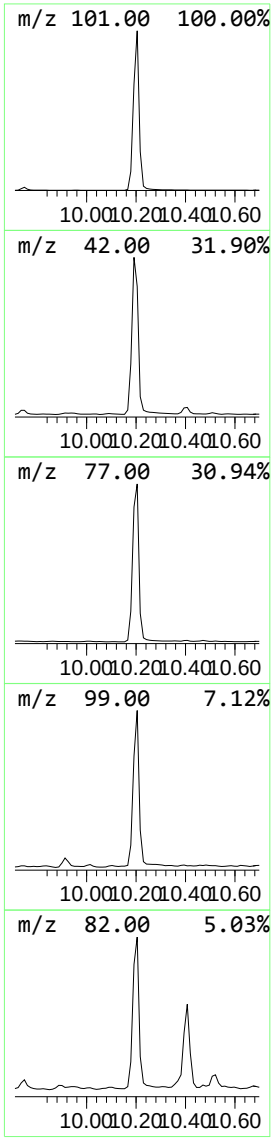
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Propane, 1-isothiocyanato- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.204	0.69 ng	315213	Naphthalene-d8	10.407

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	5
2	Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	5
3	3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	4
4	3-Aminopropionitrile	70	C3H6N2	000151-18-8	4
5	1,3,4-Thiadiazol-2-amine	101	C2H3N3S	004005-51-0	4





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Data File : BN016642.D
Acq On : 01 Oct 2021 18:57
Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

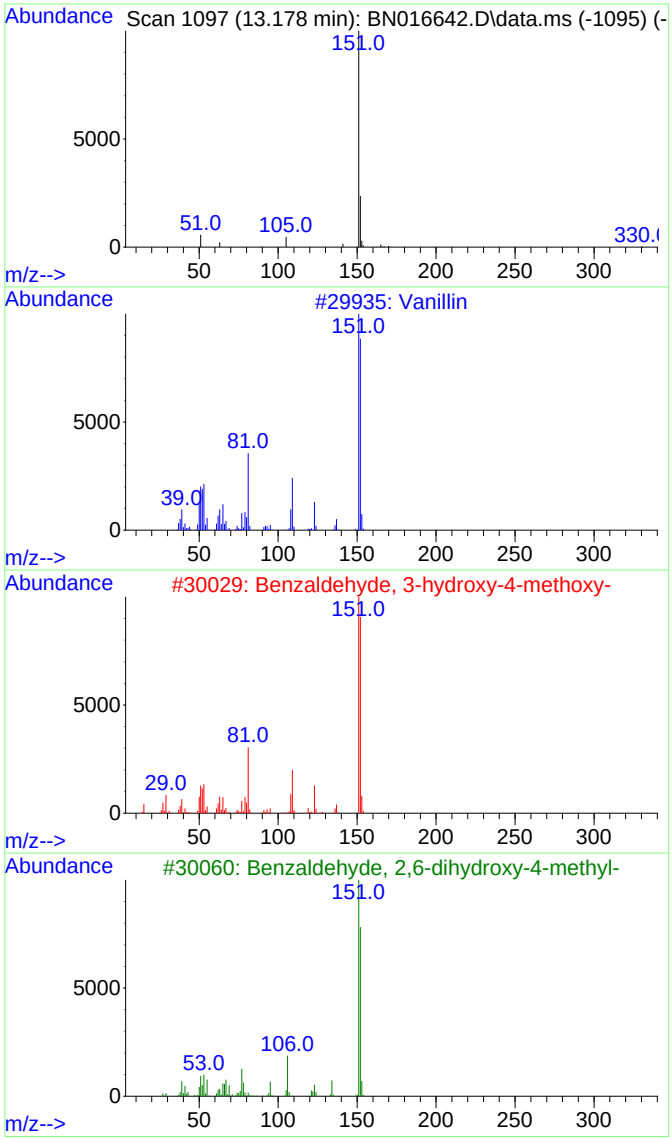
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

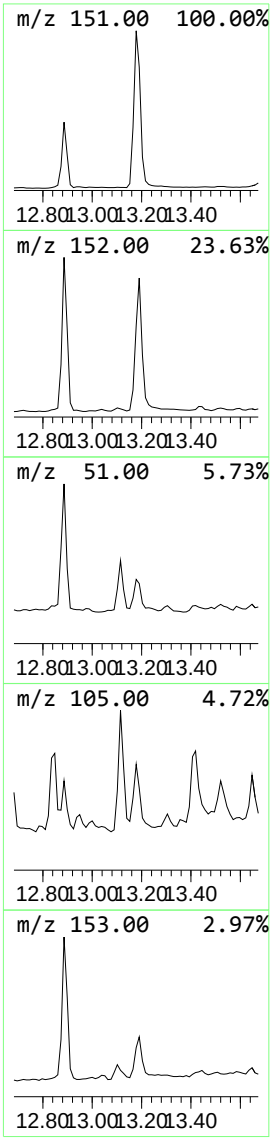
****

Peak Number 9 Vanillin Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.178	0.05 ng	23511	Acenaphthene-d10	14.269

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	5
2	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	5
3	Benzaldehyde, 2,6-dihydroxy-4-me...	152	C8H8O3	000526-37-4	5
4	3-Amino-2-methylbenzoic acid	151	C8H9NO2	052130-17-3	4
5	Furane-2-carboxaldehyde, iminose...	152	C6H8N4O	004353-50-8	4





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ALS Vial : 11 Sample Multiplier: 1

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BNA_N
ClientSampleId :
MW202D1-20210919

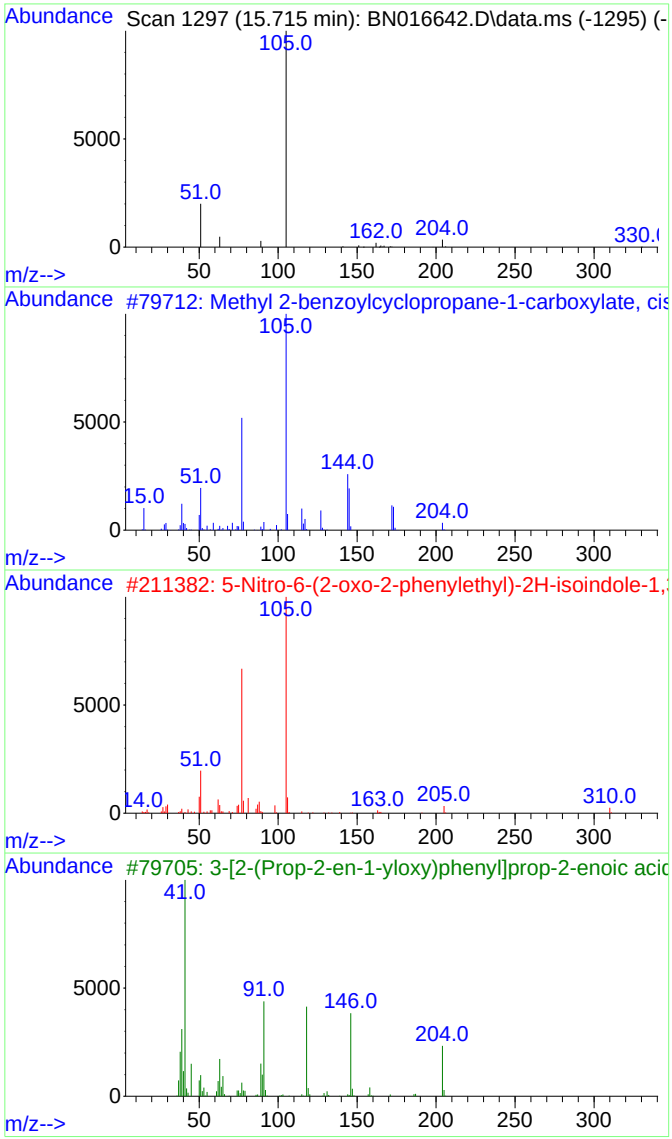
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

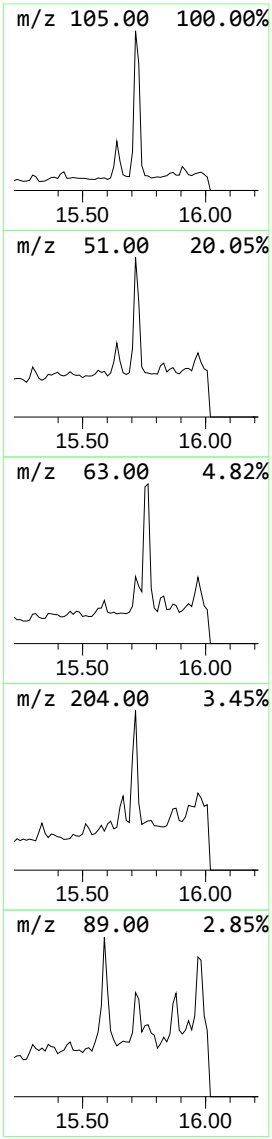
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Methyl 2-benzoylcyclopropan... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.715	0.05 ng	21845	Phenanthrene-d10	17.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 2-benzoylcyclopropane-1-c...	204	C12H12O3	006366-30-9	5
2	5-Nitro-6-(2-oxo-2-phenylethyl)-...	310	C16H10N2O5	1000489-20-6	4
3	3-[2-(Prop-2-en-1-yloxy)phenyl]p...	204	C12H12O3	091963-08-5	3
4	Phenol, 3,5-dichloro-, acetate	204	C8H6Cl2O2	061925-86-8	3
5	Phenol, 2,4-dichloro-, acetate	204	C8H6Cl2O2	006341-97-5	3





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Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

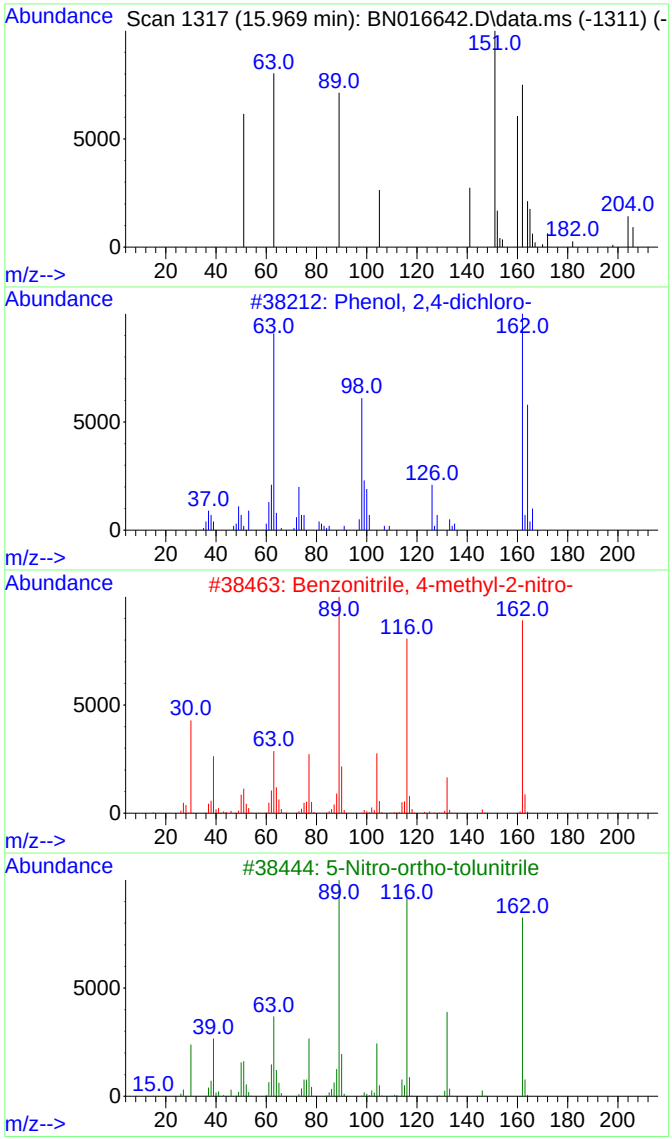
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

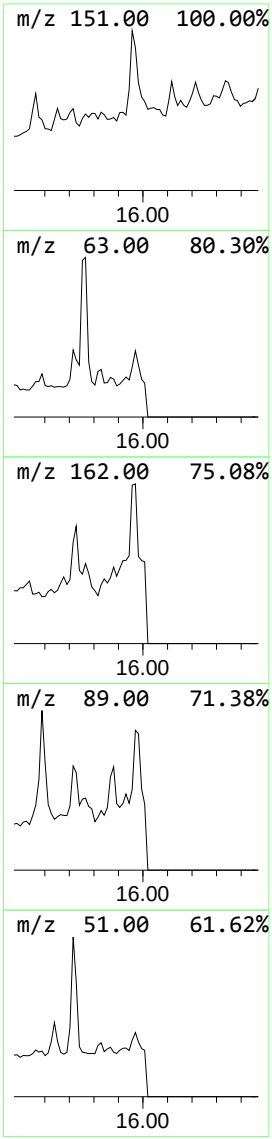
****

Peak Number 11 Phenol, 2,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.969	0.06 ng	26888	Phenanthrene-d10	17.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	9
2	Benzonitrile, 4-methyl-2-nitro-	162	C8H6N2O2	026830-95-5	7
3	5-Nitro-ortho-tolunitrile	162	C8H6N2O2	000939-83-3	7
4	6-Nitroindole	162	C8H6N2O2	004769-96-4	7
5	6-Nitro-m-tolunitrile	162	C8H6N2O2	064113-86-6	7





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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

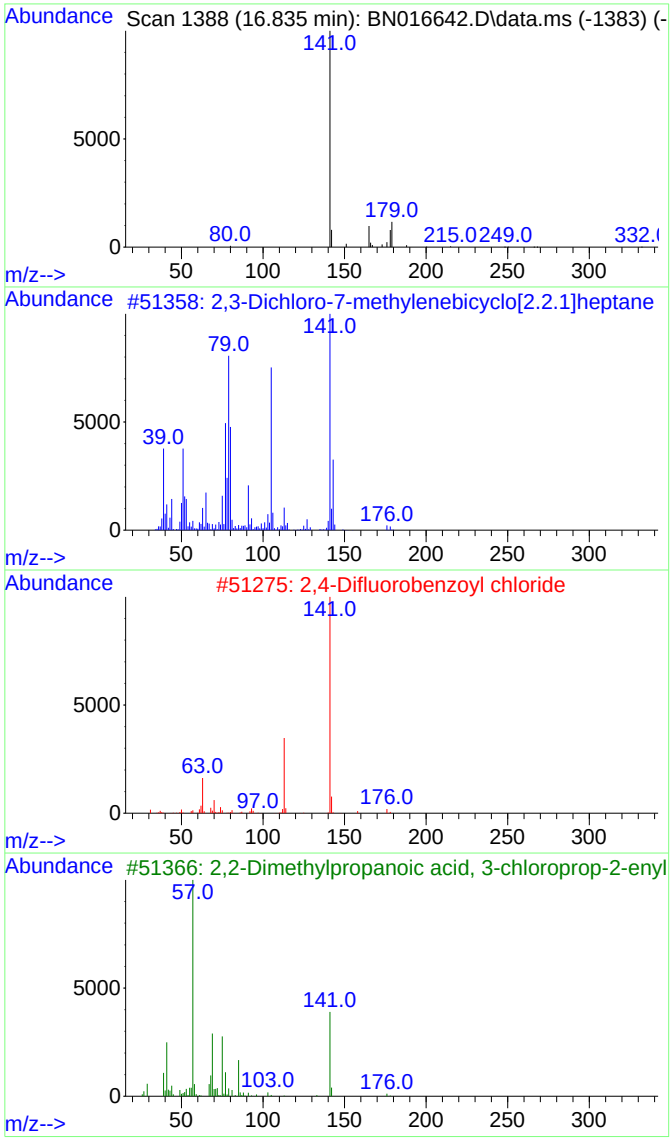
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

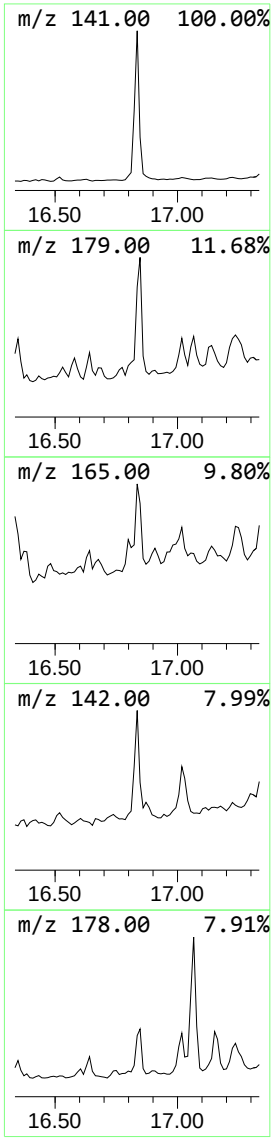
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2,3-Dichloro-7-methylenebic... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.835	0.06 ng	25867	Phenanthrene-d10	17.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dichloro-7-methylenebicyclo[...	176	C8H10Cl2	1000210-21-0	7
2	2,4-Difluorobenzoyl chloride	176	C7H3ClF2O	072482-64-5	5
3	2,2-Dimethylpropanoic acid, 3-ch...	176	C8H13ClO2	1000299-33-9	5
4	2-Chloromethyl-5-methylpyridine	141	C7H8ClN	1000241-93-8	4
5	(4-Chloro-1H-pyrazol-1-yl)aceton...	141	C5H4ClN3	113336-23-5	4





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Operator : CG/JU
Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

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

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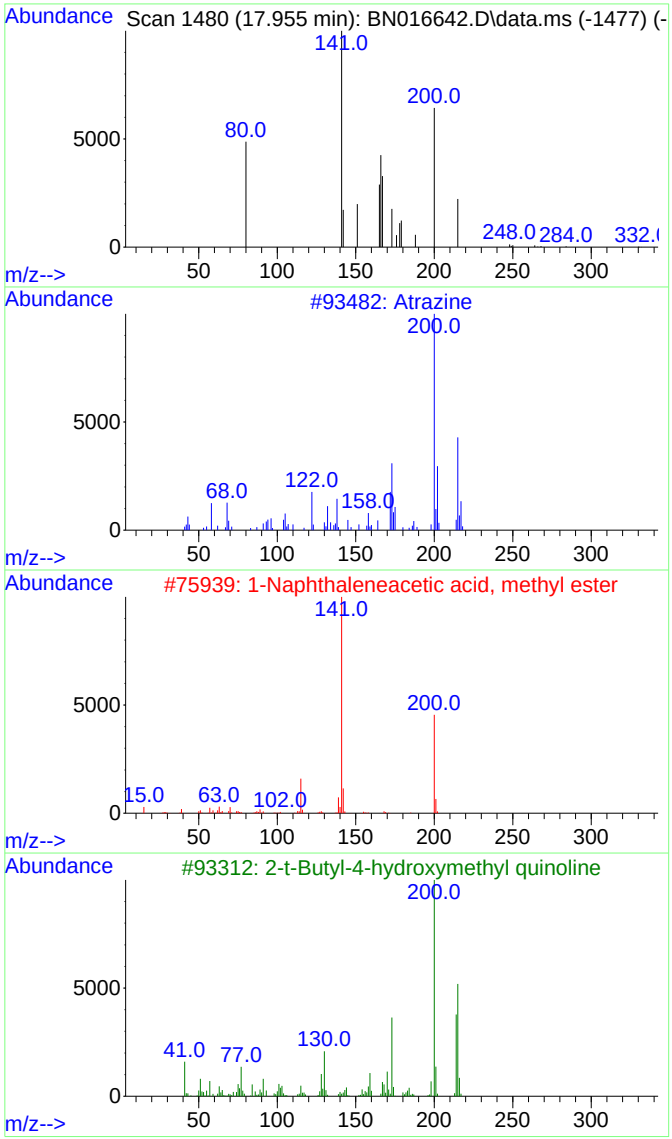
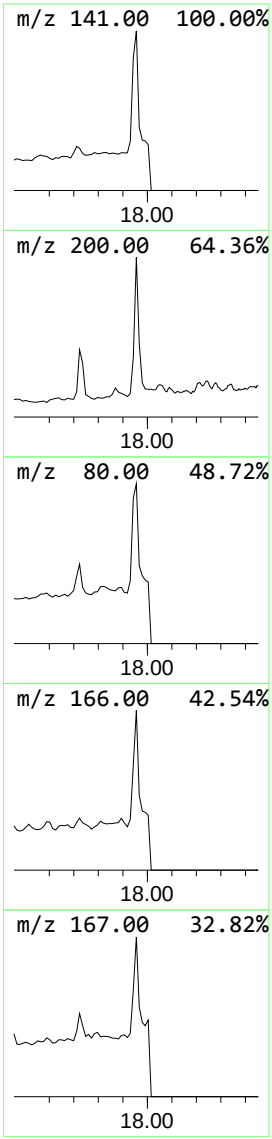
Peak Number 13 Atrazine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.955	0.06 ng	24932	Phenanthrene-d10	17.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Atrazine	215	C8H14ClN5	001912-24-9	9
2	1-Naphthaleneacetic acid, methyl...	200	C13H12O2	002876-78-0	9
3	2-t-Butyl-4-hydroxymethyl quinoline	215	C14H17NO	1000255-75-0	9
4	Hydrazinecarboxamide, 2-cyclopen...	141	C6H11N3O	005459-00-7	9
5	4-Chlorophenoxyacetic acid, meth...	200	C9H9ClO3	1000362-94-5	9





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Sample : M3833-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

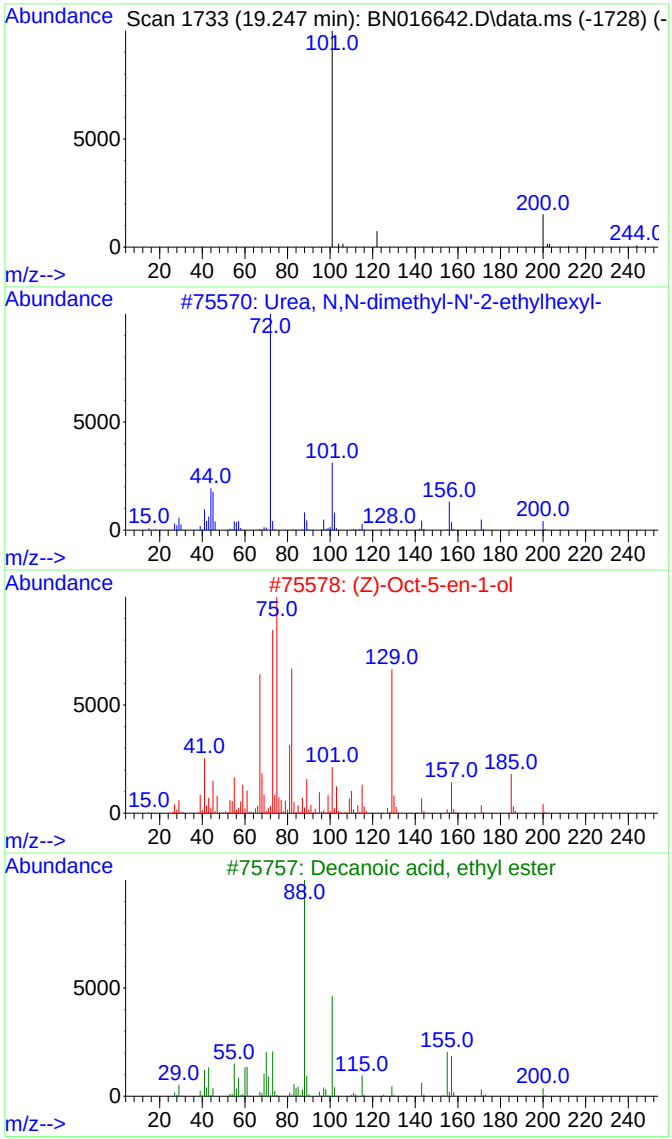
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

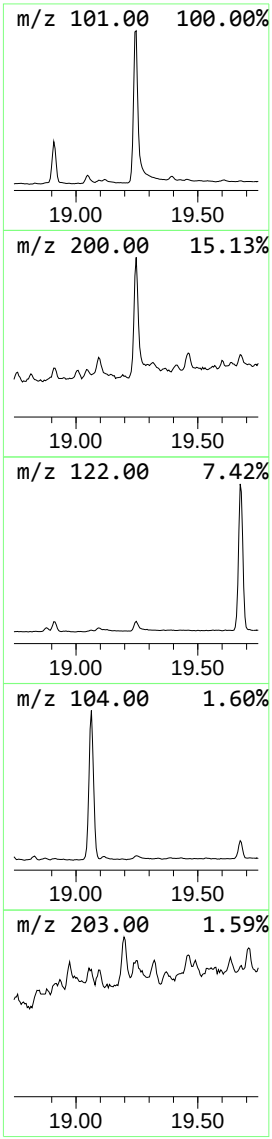
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Urea, N,N-dimethyl-N'-2-eth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.247	0.04 ng	22049	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Urea, N,N-dimethyl-N'-2-ethylhexyl-	200	C11H24N2O	1010419-88-4	4
2	(Z)-Oct-5-en-1-ol	200	C11H24OSi	1010498-20-3	4
3	Decanoic acid, ethyl ester	200	C12H24O2	000110-38-3	4
4	Octan-2-yl (E)-2-methylbut-2-enoate	212	C13H24O2	1000373-76-0	3
5	Silane, diethylmethyl-7-octenyl-	212	C13H28Si	062185-52-8	3





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Misc :
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Instrument :
BNA_N
ClientSampleId :
MW202D1-20210919

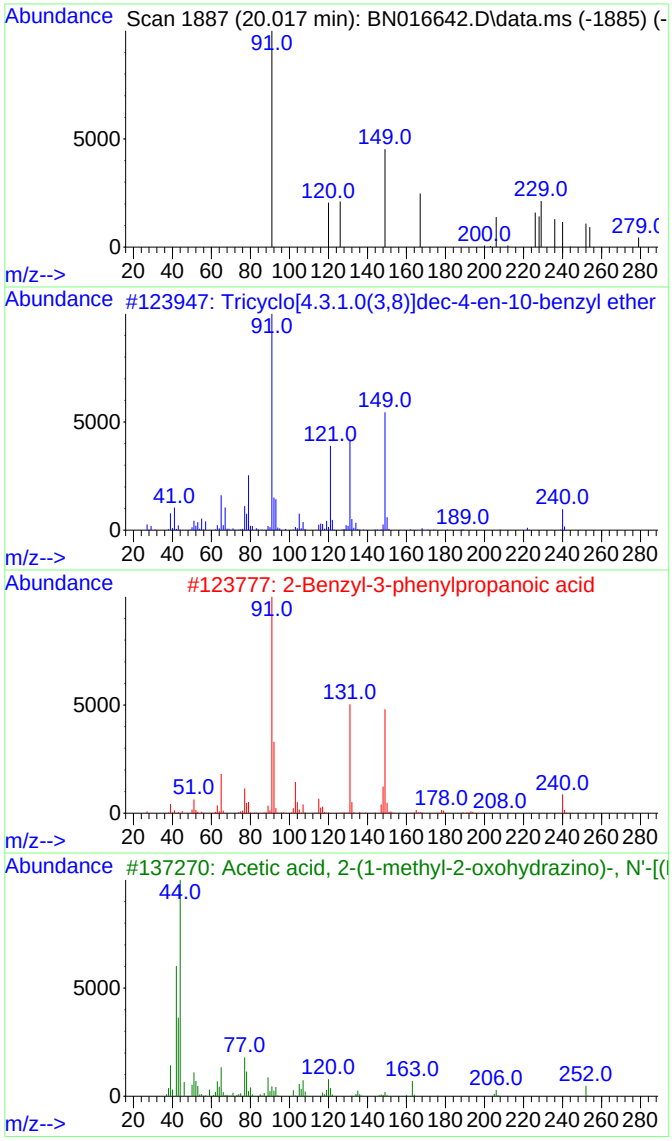
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

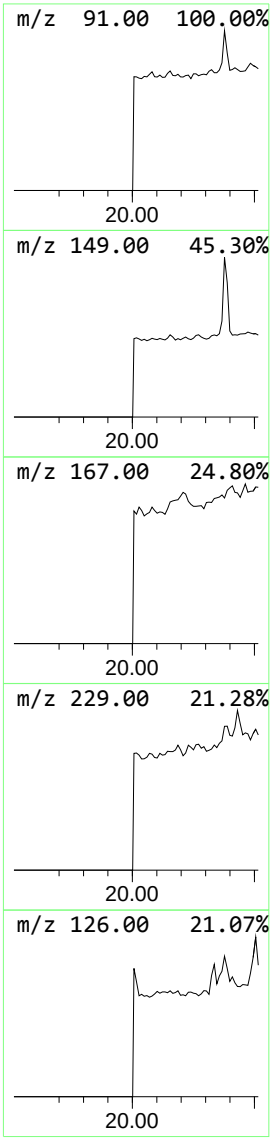
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Tricyclo[4.3.1.0(3,8)]dec-4... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.017	0.06 ng	28484	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[4.3.1.0(3,8)]dec-4-en-1...	240	C17H20O	1000197-67-3	9
2	2-Benzyl-3-phenylpropanoic acid	240	C16H16O2	000618-68-8	7
3	Acetic acid, 2-(1-methyl-2-oxohy...	252	C10H12N4O4	1000327-59-6	5
4	Bicyclo[2.2.1]hept-5-ene-2-aceti...	236	C13H16O4	077828-47-8	5
5	2,2,5,5-Tetramethyl-4-(p-fluorop...	236	C13H17FN2O	090277-88-6	5





Tentatively Identified Compound (LSC) summary

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1,4-Butanediamine	3.502	0.2	ng	42930	1	7.643	106635	0.4
1-Propanol, 2-m...	3.825	0.2	ng	48271	1	7.643	106635	0.4
Butane	4.821	0.5	ng	134796	1	7.643	106635	0.4
Acetic acid, me...	6.656	0.1	ng	19353	1	7.643	106635	0.4
Hydroxylamine, ...	7.265	0.1	ng	25636	1	7.643	106635	0.4
2,3-Octanedione	7.864	0.1	ng	19790	1	7.643	106635	0.4
5-Hepten-2-one	8.177	0.0	ng	11213	1	7.643	106635	0.4
Propane, 1-isot...	10.204	0.7	ng	315213	2	10.407	182736	0.4
Vanillin	13.178	0.0	ng	23511	3	14.269	202400	0.4
Methyl 2-benzoy...	15.715	0.1	ng	21845	4	17.030	167423	0.4
Phenol, 2,4-dic...	15.969	0.1	ng	26888	4	17.030	167423	0.4
2,3-Dichloro-7-...	16.835	0.1	ng	25867	4	17.030	167423	0.4
Atrazine	17.955	0.1	ng	24932	4	17.030	167423	0.4
Urea, N,N-dimet...	19.247	0.0	ng	22049	5	21.228	199965	0.4
Tricyclo[4.3.1....	20.017	0.1	ng	28484	5	21.228	199965	0.4