LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : S-829-J-SO-3-3.5-092221

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: BN016685.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.512	44	47	58	rVB	7445	24478	8.41%	0.743%
2	3.825	78	81	88	rVB	5015	11062	3.80%	0.336%
3	4.821	185	189	202	rVB	33209	70543	24.23%	2.141%
4	5.042	210	213	220	rBV	4651	11561	3.97%	0.351%
5	5.273	234	238	246	rVB	16404	35581	12.22%	1.080%
6	6.647	384	387	391	rBV	3129	5374	1.85%	0.163%
7	6.831	403	407	421	rVB	17394	39364	13.52%	1.194%
8	7.264	450	454	459	rVB2	3473	7845	2.69%	0.238%
9	7.643	490	495	500	rBV	47289	94584	32.49%	2.870%
10	7.707	500	502	508	rVB	1836	3125	1.07%	0.095%
11	7.938	523	527	532	rVB	10097	17852	6.13%	0.542%
12	8.306	565	567	569	rVB	5268	5207	1.79%	0.158%
13	8.417	574	576	580	rVB3	1838	3743	1.29%	0.114%
14	8.785	601	605	612	rBV	22990	45595	15.66%	1.384%
15	8.949	616	618	622	rVB	3094	4819	1.66%	0.146%
16	10.191	713	716	723	rBV	33012	57720	19.83%	1.752%
17	10.407	729	733	736	rBV	100898	173562	59.62%	5.267%
18	10.508	739	741	745	rVB2	4127	7075	2.43%	0.215%
19	12.003	921	931	943	rBV	39828	65572	22.52%	1.990%
20	12.886	1071	1074	1078	rBV2	75763	128290	44.07%	3.893%
21	13.178	1094	1097	1102	rBV2	7687	14803	5.08%	0.449%
22	14.269	1180	1183	1186	rBV	137276	197046	67.69%	5.979%
23	14.332	1186	1188	1192	rVV	14882	22737	7.81%	0.690%
24	14.434	1193	1196	1200	rVB	5059	7747	2.66%	0.235%
25	14.523	1200	1203	1207	rBV3	1908	4332	1.49%	0.131%
26	15.322	1263	1266	1269	rBV	3928	6174	2.12%	0.187%
27	15.639	1288	1291	1295	rBV2	2444	5668	1.95%	0.172%
28	15.715	1295	1297	1299	rVV	3664	6251	2.15%	0.190%
29	15.766	1299	1301	1306	rVB2	10622	17068	5.86%	0.518%
30	16.007	1317	1320	1322	rVB3	4692	7099	2.44%	0.215%
31	16.835	1385	1388	1391	rVB	2937	4503	1.55%	0.137%
32	17.018	1400	1403	1405	rBV	105272	160461	55.12%	4.869%
33	17.066	1405	1407	1410	rVV	16639	26578	9.13%	0.807%
34	17.152	1411	1414	1418	rVB	5221	8737	3.00%	0.265%
35	17.711	1456	1460	1465	rBV3	1389	3707	1.27%	0.112%
36	17.943	1477	1479	1483	rBV	3426	5393	1.85%	0.164%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : S-829-J-SO-3-3.5-092221

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	19.063	1688	1696	1699	rBV	98665	131962	45.33%	4.004%
38	19.093	1699	1702	1710	rVB	60932	78517	26.97%	2.383%
39	19.242	1729	1732	1745	rVB3	3450	7181	2.47%	0.218%
40	19.460	1766	1776	1787	rBV	51787	75332	25.88%	2.286%
41	19.679	1814	1820	1836	rVB	84006	108545	37.29%	3.294%
42	20.006	1884	1886	1888	rBV	4952	6246	2.15%	0.190%
43	20.378	1919	1921	1923	rBV	2727	4446	1.53%	0.135%
44	20.474	1923	1930	1940	rVB3	10159	49085	16.86%	1.489%
45	20.920	1970	1972	1973	rBV	2596	3067	1.05%	0.093%
46	20.973	1974	1977	1980	rVV	34579	60193	20.68%	1.827%
47	21.090	1980	1988	1992	rVV3	25117	133174	45.75%	4.041%
48	21.164	1993	1995	1998	rVV	32988	64187	22.05%	1.948%
49	21.228	1998	2001	2003	rVV2	163248	291118	100.00%	8.834%
50	21.270	2003	2005	2011	rVB	45017	66211	22.74%	2.009%
51	21.674	2040	2043	2047	rBV2	4730	8351	2.87%	0.253%
52	21.875	2057	2062	2065	rBV2	65204	106592	36.61%	3.235%
53	22.003	2071	2074	2077	rVB	35803	50524	17.36%	1.533%
54	22.109	2081	2084	2088	rVB3	7807	12294	4.22%	0.373%
55	22.226	2093	2095	2099	rVV	6037	19609	6.74%	0.595%
56	22.321	2101	2104	2106	rVV	9378	22285	7.65%	0.676%
57	22.396	2106	2111	2112	rVB	26893	51081	17.55%	1.550%
58	22.807	2219	2229	2235	rBV	32863	61507	21.13%	1.866%
59	22.848	2237	2241	2251	rVB2	27752	43537	14.96%	1.321%
60	22.978	2273	2279	2303	rVB2	5070	10167	3.49%	0.309%
61	23.286	2361	2369	2385	rVB	11491	21023	7.22%	0.638%
62	23.382	2387	2397	2405	rBV	23136	46351	15.92%	1.407%
63	23.481	2415	2426	2464	rVB	111127	229041	78.68%	6.950%
64	24.084	2593	2602	2618	rVB	9465	18571	6.38%	0.564%
65	24.186	2622	2632	2652	rBV	10379	31633	10.87%	0.960%
66	24.857	2818	2828	2855	rVB	8044	19800	6.80%	0.601%
67	25.466	2996	3006	3015	rBV4	1185	3251	1.12%	0.099%
68	25.768	3076	3094	3121	rBV3	21483	75679	26.00%	2.296%
69	26.456	3278	3295	3333	rVB	12457	40584	13.94%	1.232%
70	26.825	3390	3403	3436	rVB3	2937	10633	3.65%	0.323%
71	27.483	3577	3595	3626	rVB3	6233	22415	7.70%	0.680%

Sum of corrected areas: 3295448

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Acq On : 02 Oct 2021 23:45
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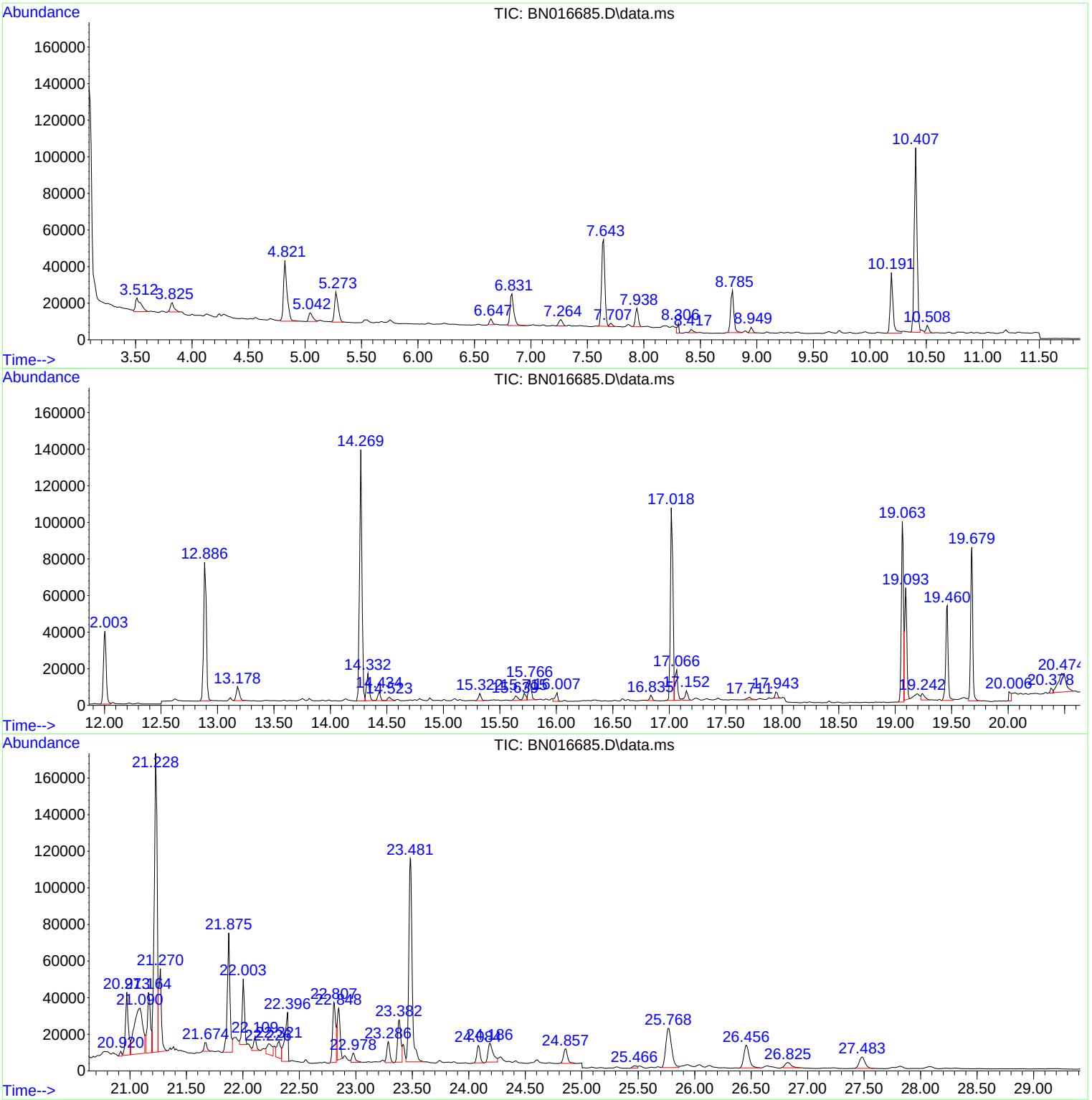
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

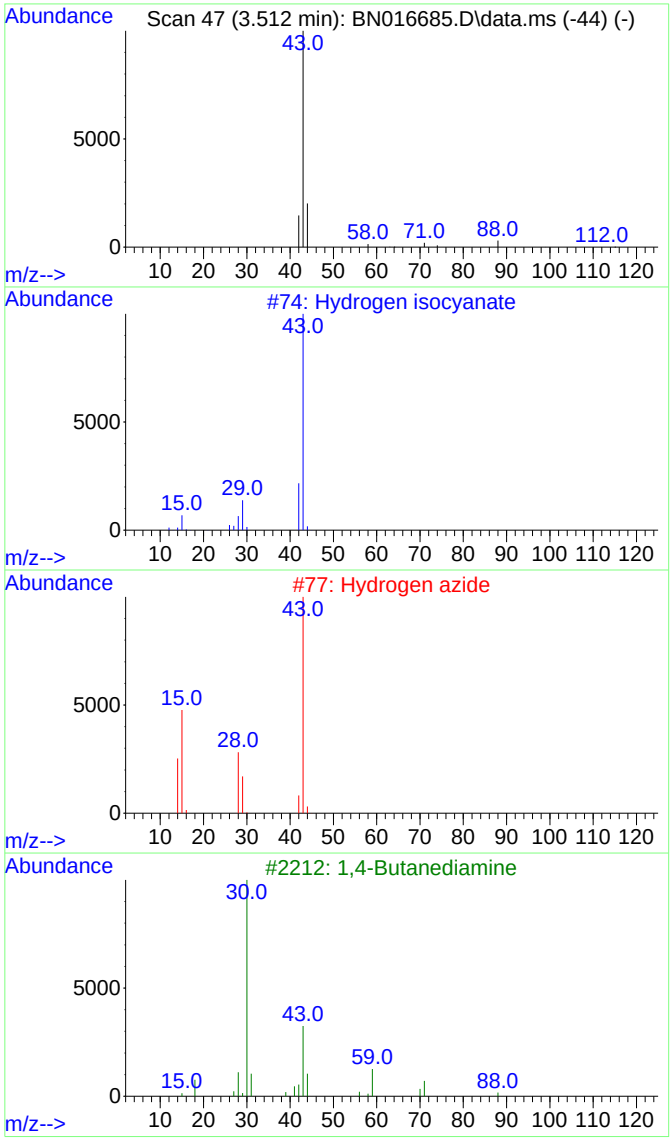
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

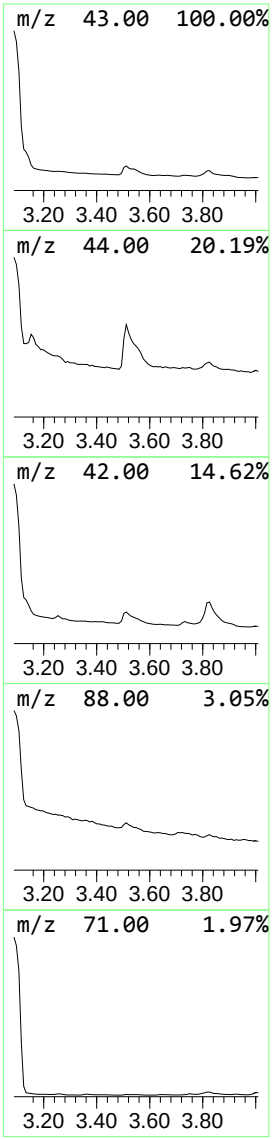
****

Peak Number 1 Hydrogen isocyanate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.512	0.10 ng	24478	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrogen isocyanate	43	CHNO	000075-13-8	3
2	Hydrogen azide	43	HN3	007782-79-8	3
3	1,4-Butanediamine	88	C4H12N2	000110-60-1	3
4	Isobutane	58	C4H10	000075-28-5	2
5	Guanidine, methyl-	73	C2H7N3	000471-29-4	2





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

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

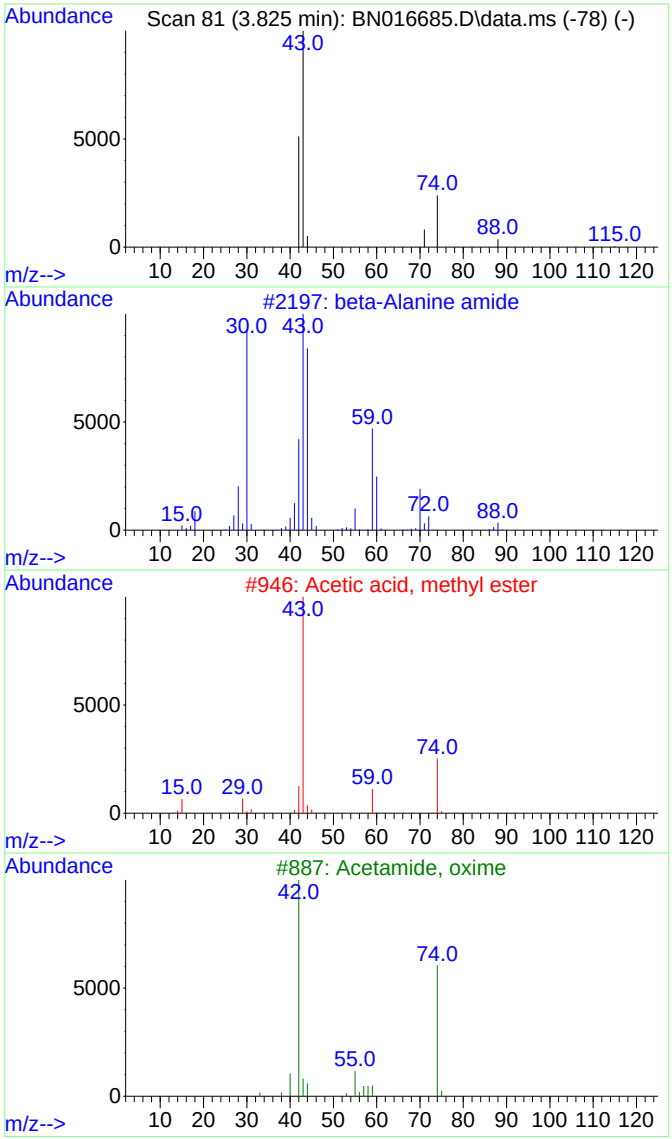
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

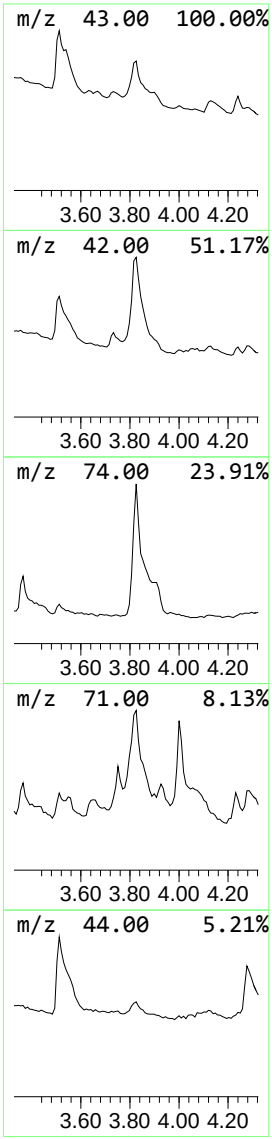
****

Peak Number 2 beta-Alanine amide Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.825	0.05 ng	11062	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	beta-Alanine amide	88	C3H8N2O	004726-85-6	5
2	Acetic acid, methyl ester	74	C3H6O2	000079-20-9	5
3	Acetamide, oxime	74	C2H6N2O	022059-22-9	4
4	O-Methylisourea	74	C2H6N2O	002440-60-0	3
5	1-Propyne, 3-chloro-	74	C3H3Cl	000624-65-7	3





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Data File : BN016685.D
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Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

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

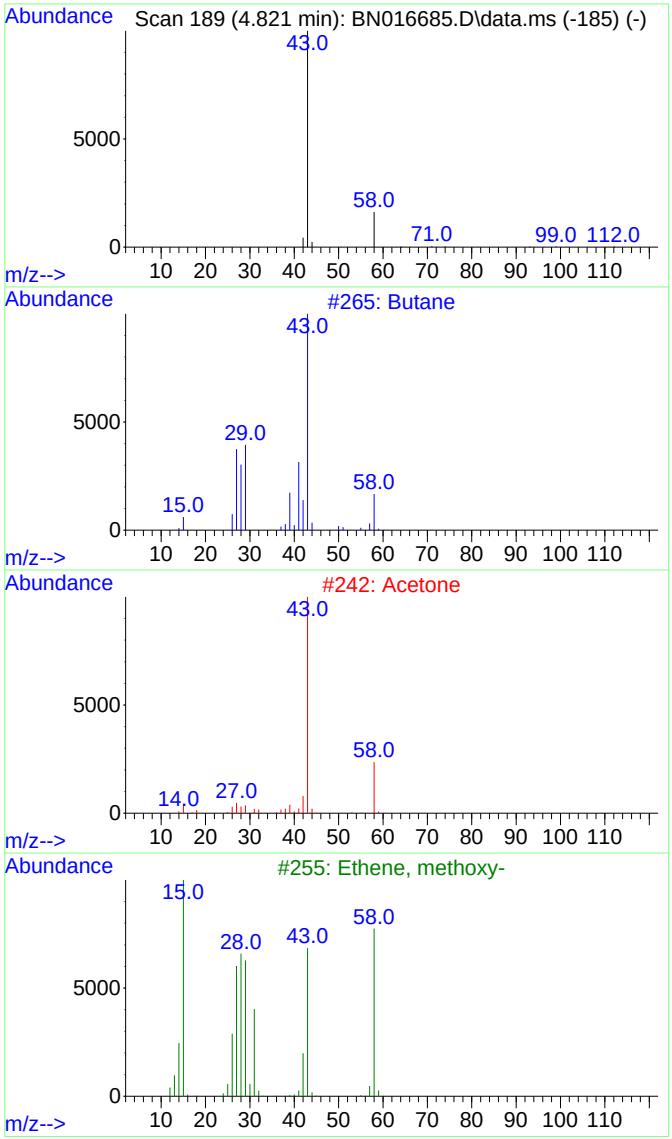
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

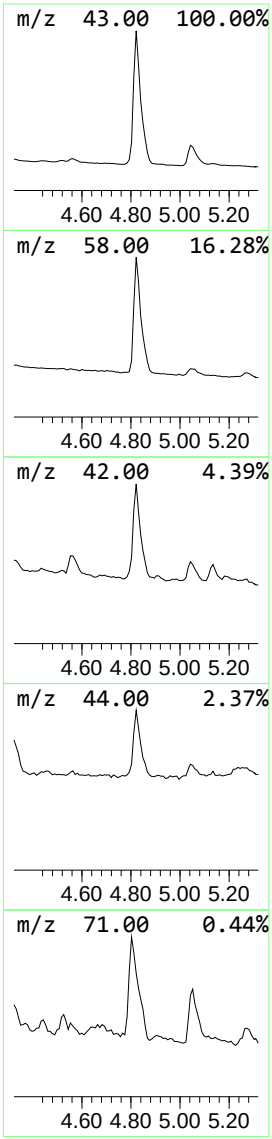
****

Peak Number 3 Butane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.821	0.30 ng	70543	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\
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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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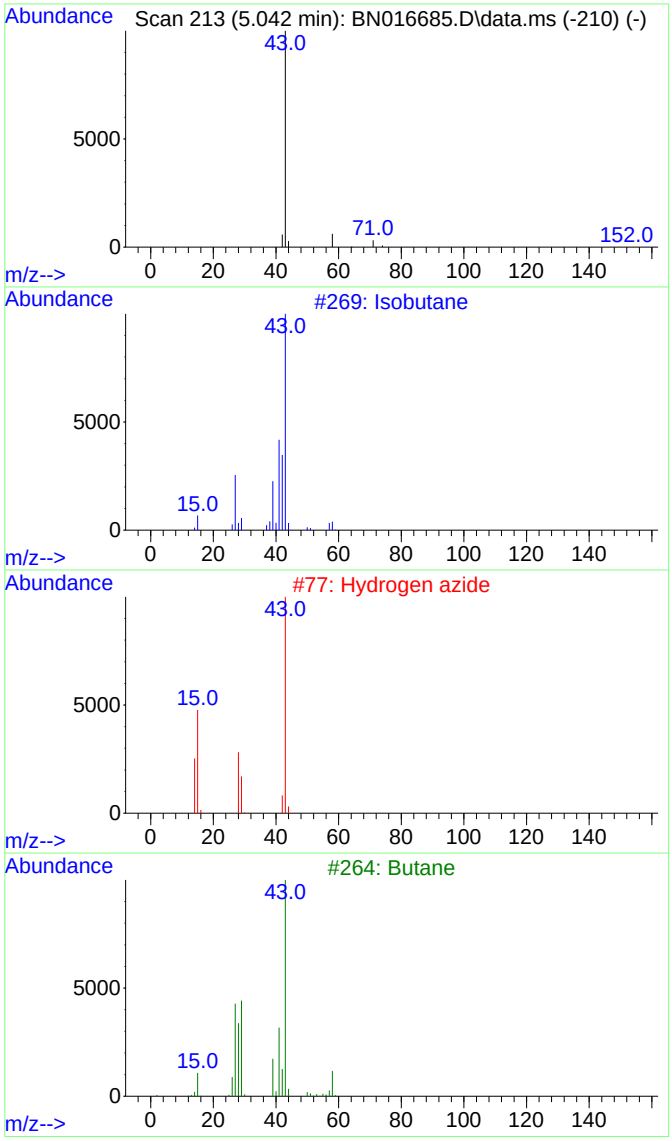
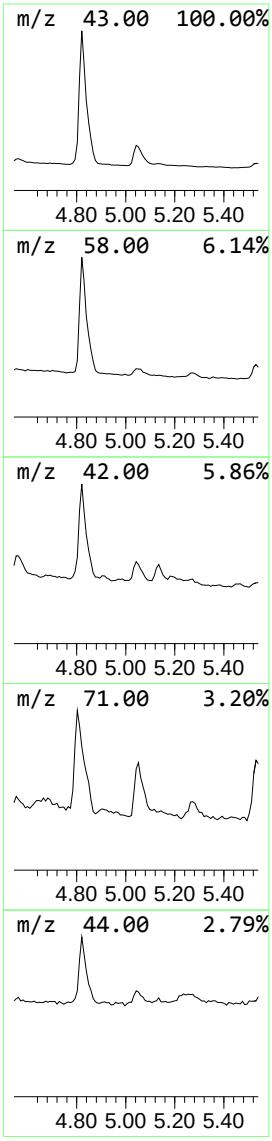
****

Peak Number 4 Isobutane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.042	0.05 ng	11561	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutane	58	C4H10	000075-28-5	4
2	Hydrogen azide	43	HN3	007782-79-8	4
3	Butane	58	C4H10	000106-97-8	3
4	Hydrogen isocyanate	43	CHNO	000075-13-8	2
5	Methoxyacetonitrile	71	C3H5NO	001738-36-9	2





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Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
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Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

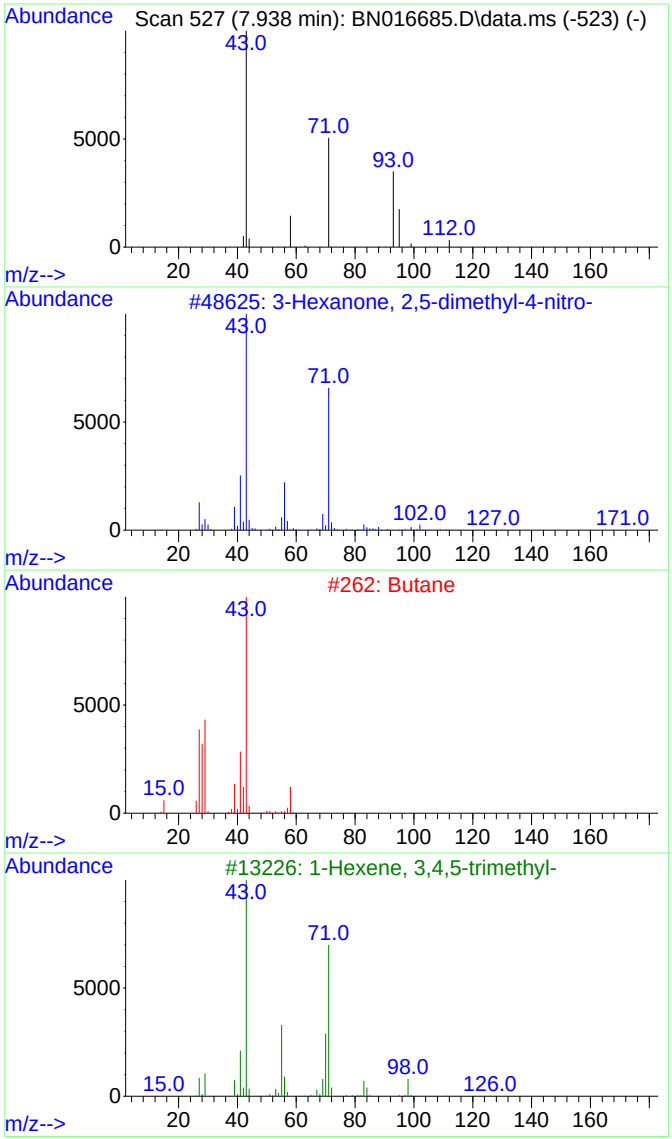
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

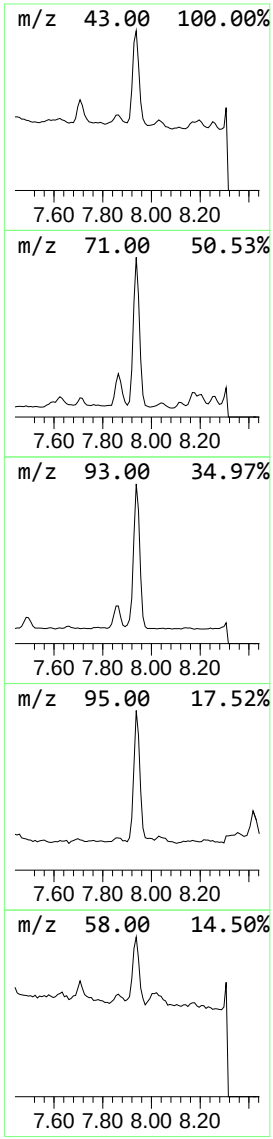
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 3-Hexanone, 2,5-dimethyl-4-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.938	0.08 ng	17852	1,4-Dichlorobenzene-d4	7.643

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone, 2,5-dimethyl-4-nitro-	173	C8H15NO3	059906-54-6	9
2	Butane	58	C4H10	000106-97-8	4
3	1-Hexene, 3,4,5-trimethyl-	126	C9H18	056728-10-0	4
4	2-Furancarboxylic acid, 2-tetrah...	196	C10H12O4	004623-04-5	4
5	Cyclopropanemethanol, .alpha.-bu...	128	C8H16O	004379-16-2	4





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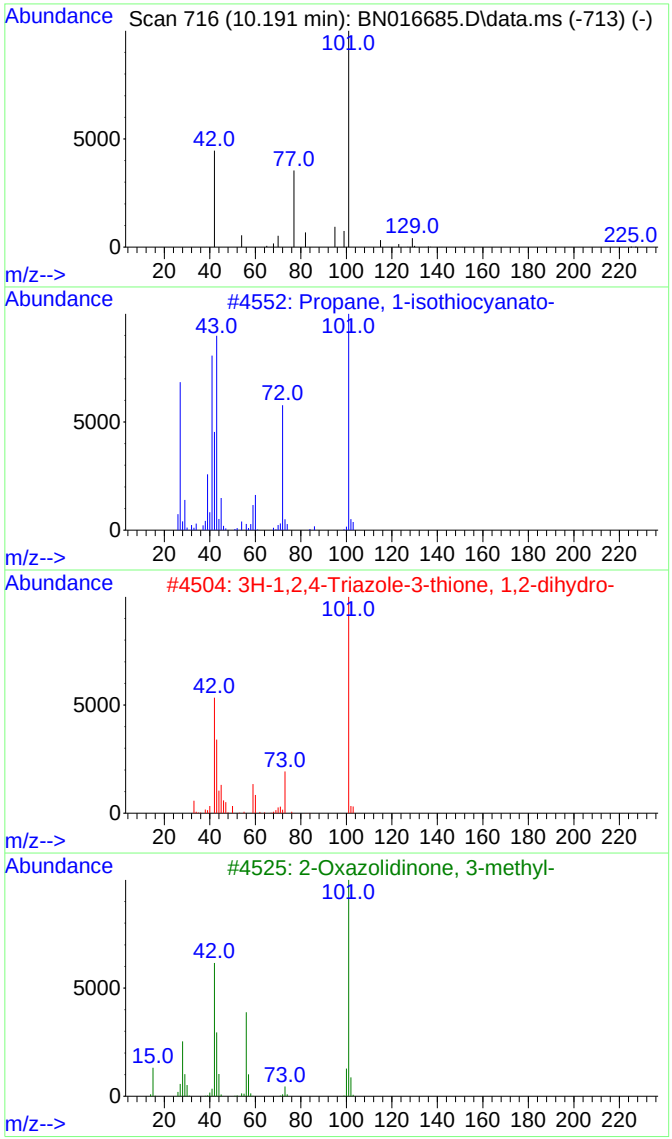
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

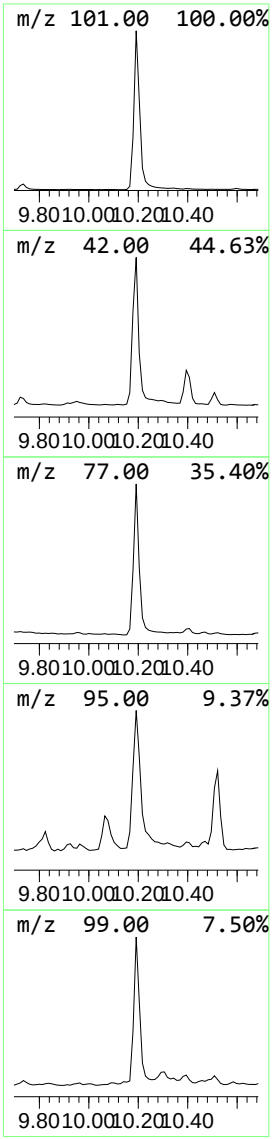
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Propane, 1-isothiocyanato- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.191	0.13 ng	57720	Naphthalene-d8	10.407

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	5
2	3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	5
3	2-Oxazolidinone, 3-methyl-	101	C4H7NO2	019836-78-3	5
4	Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	5
5	3-Aminopropionitrile	70	C3H6N2	000151-18-8	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

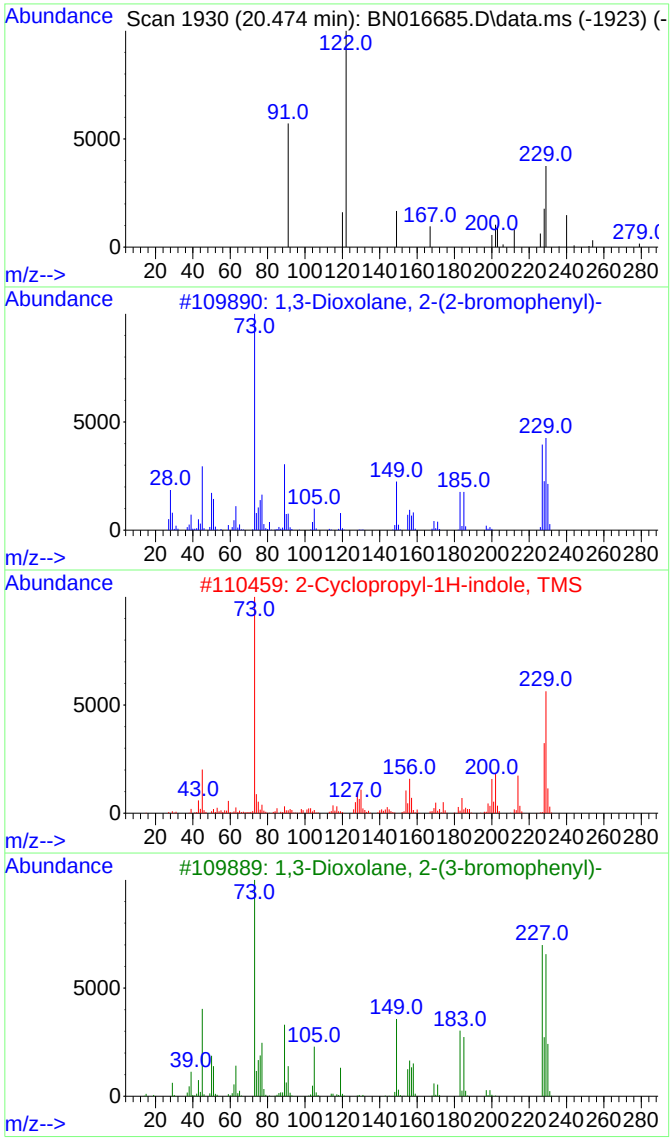
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

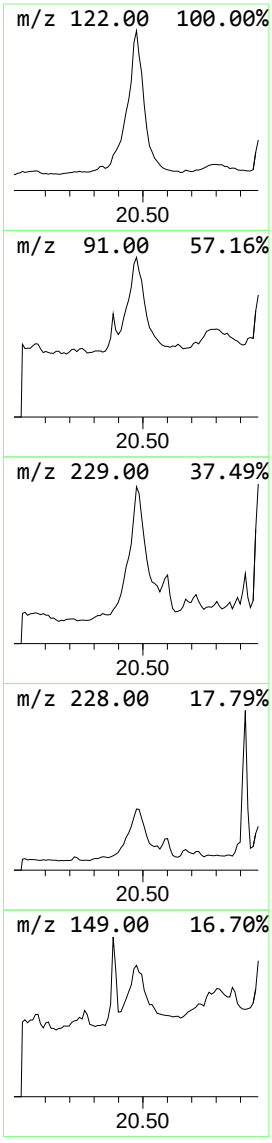
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,3-Dioxolane, 2-(2-bromoph... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.474	0.07 ng	49085	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(2-bromophenyl)-	228	C9H9BrO2	034824-58-3	9
2	2-Cyclopropyl-1H-indole, TMS	229	C14H19NSi	1000494-09-0	9
3	1,3-Dioxolane, 2-(3-bromophenyl)-	228	C9H9BrO2	017789-14-9	7
4	2-Chloro-6-phenoxybenzonitrile	229	C13H8ClNO	1000340-54-1	7
5	N-(4,6-Dimethyl-2-pyridinyl)-2-p...	240	C15H16N2O	127722-57-0	5





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Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

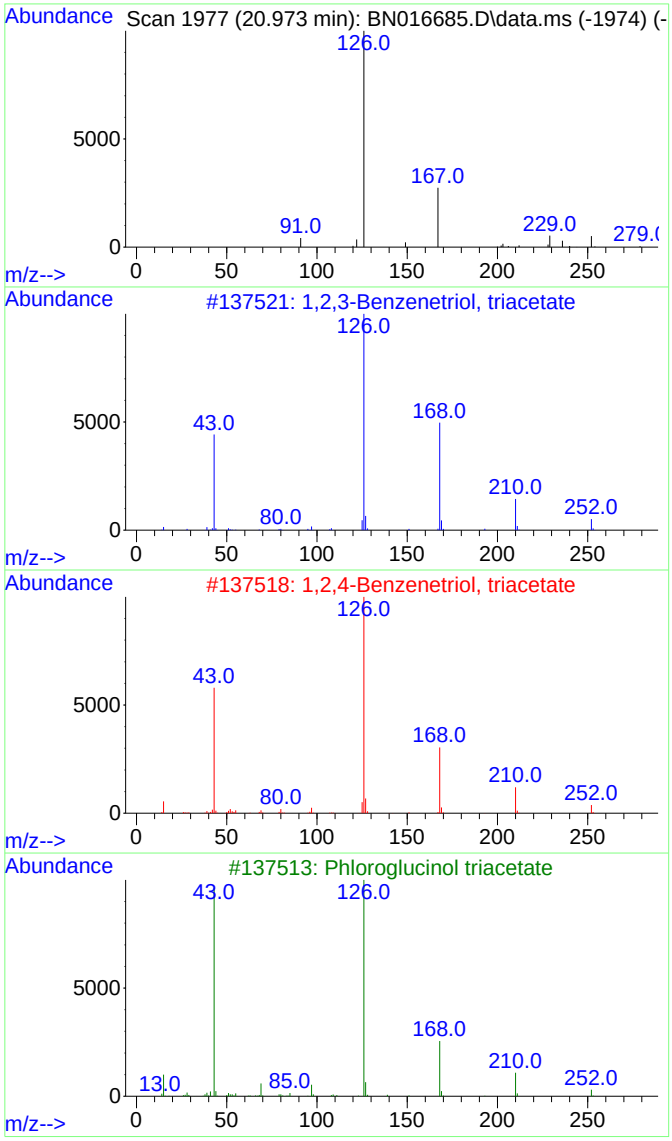
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

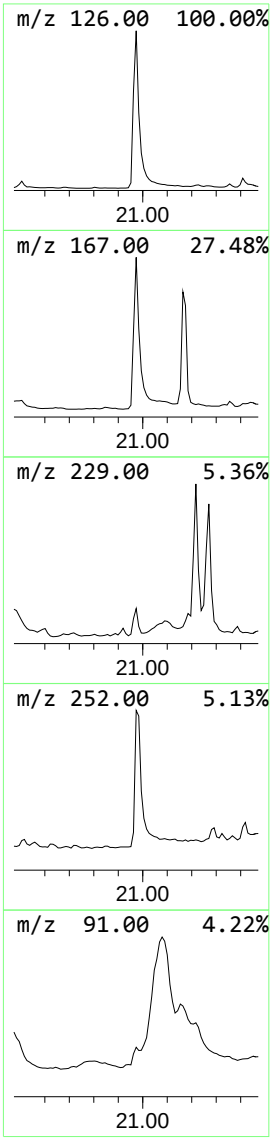
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,2,3-Benzenetriol, triacetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.973	0.08 ng	60193	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3-Benzenetriol, triacetate	252	C12H12O6	000525-52-0	4
2	1,2,4-Benzenetriol, triacetate	252	C12H12O6	000613-03-6	4
3	Phloroglucinol triacetate	252	C12H12O6	002999-40-8	4
4	(6S,12aR,E)-2,2-Dimethyltetrahyd...	236	C14H24N2O	155975-43-2	4
5	N-[[6-Cyclooctylaminohexyl]aziri...	252	C16H32N2	1010256-12-1	4





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Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

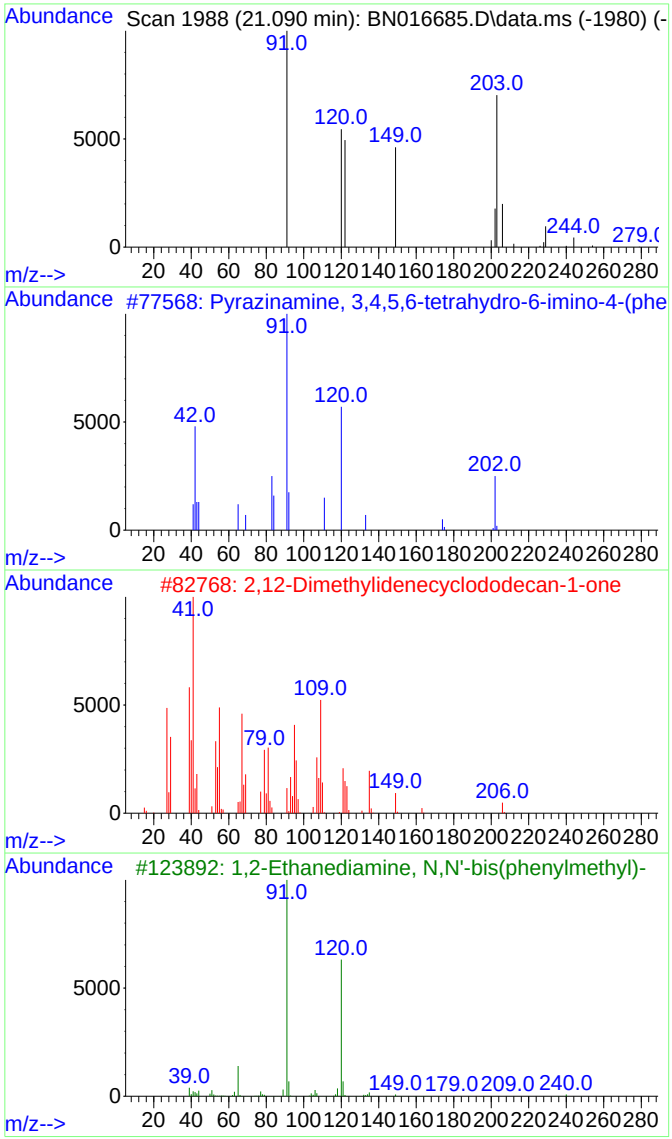
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

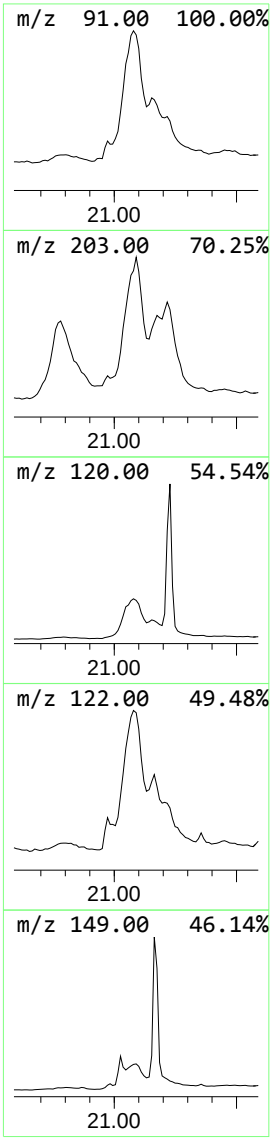
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Pyrazinamine, 3,4,5,6-tetra... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.090	0.18 ng	133174	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrazinamine, 3,4,5,6-tetrahydro...	202	C11H14N4	035975-20-3	9
2	2,12-Dimethylidenecyclododecan-1...	206	C14H22O	069915-61-3	7
3	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	7
4	Formamide, N-(2,4-dimethylphenyl)-	149	C9H11NO	060397-77-5	5
5	1,5-Benzoxazepine, 2,3,4,5-tetra...	149	C9H11NO	007160-97-6	5





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Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

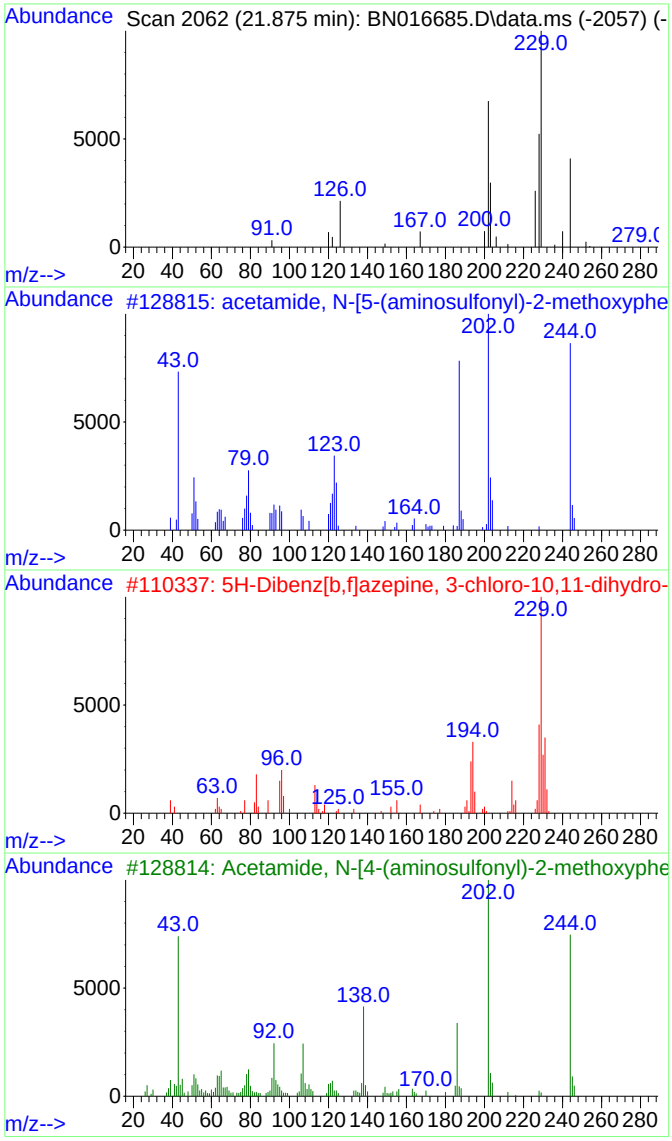
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

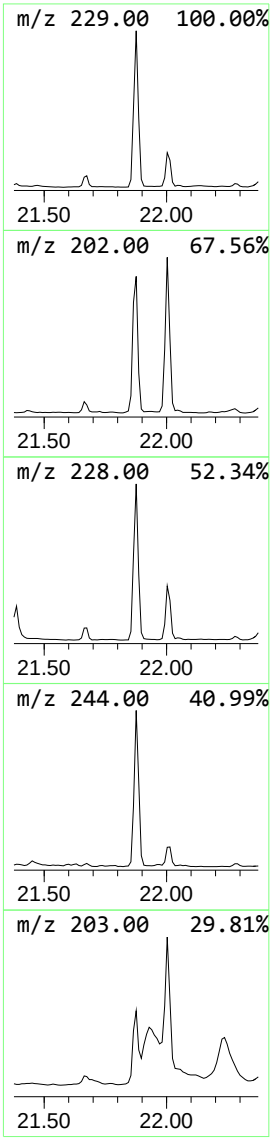
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 acetamide, N-[5-(aminosulfo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.875	0.15 ng	106592	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	acetamide, N-[5-(aminosulfonyl)-...	244	C9H12N2O4S	1000401-26-6	9
2	5H-Dibenz[b,f]azepine, 3-chloro-...	229	C14H12ClN	032943-25-2	9
3	Acetamide, N-[4-(aminosulfonyl)-...	244	C9H12N2O4S	1000107-07-5	9
4	2-Propenoic acid, 2-cyano-3-[4-(...	244	C14H16N2O2	1000115-72-8	9
5	2-Hydroxy-3,4-tetramethylene-6-m...	229	C14H15NO2	004562-57-6	9





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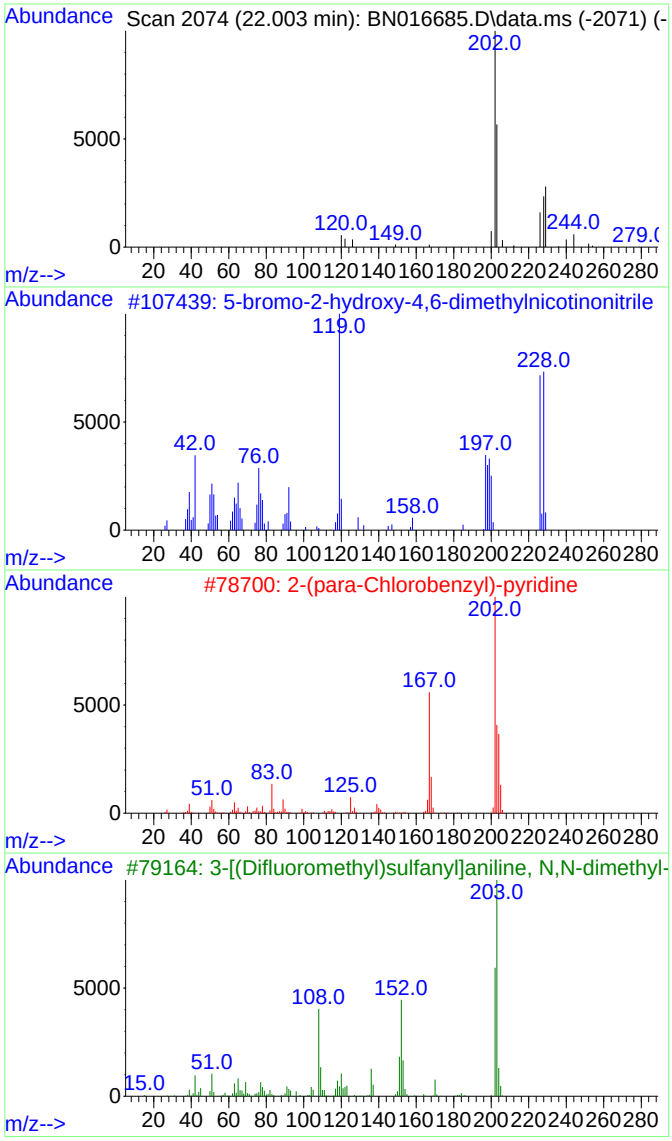
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

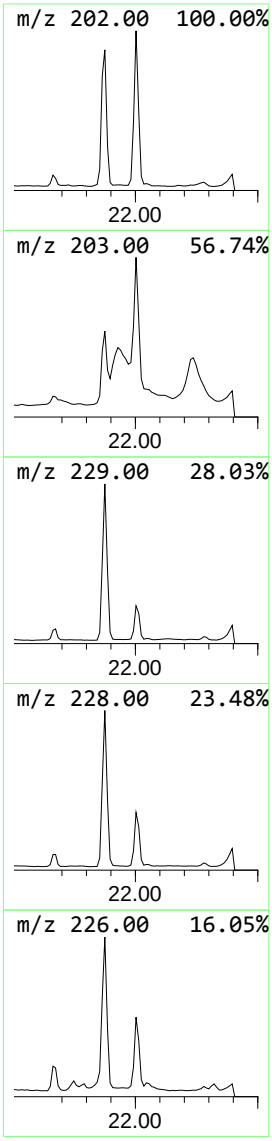
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 5-bromo-2-hydroxy-4,6-dimet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.003	0.07 ng	50524	Chrysene-d12	21.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-bromo-2-hydroxy-4,6-dimethylni...	226	C8H7BrN2O	1000472-47-7	9
2	2-(para-Chlorobenzyl)-pyridine	203	C12H10ClN	004350-41-8	9
3	3-[(Difluoromethyl)sulfanyl]anil...	203	C9H11F2NS	1000505-37-1	7
4	1H-Pyrimidin-2-one, 4-[N'-(1-pyr...	229	C11H11N5O	1000304-67-8	7
5	2-Chloro-5-(methylthio)benzoic acid	202	C8H7ClO2S	051546-12-4	7





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Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

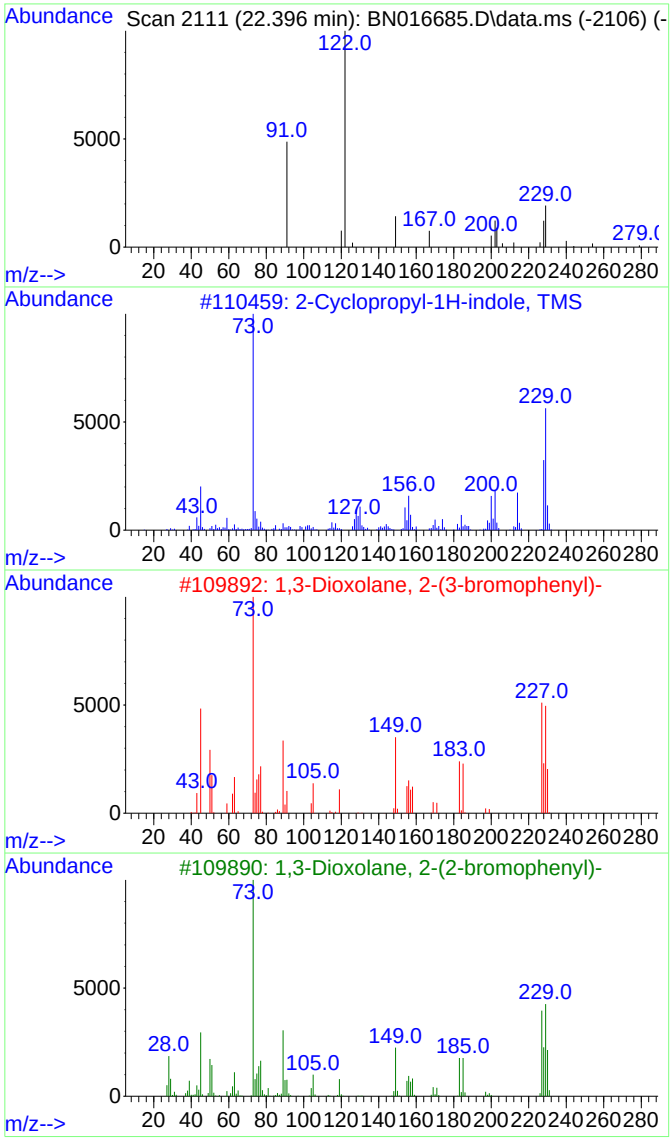
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

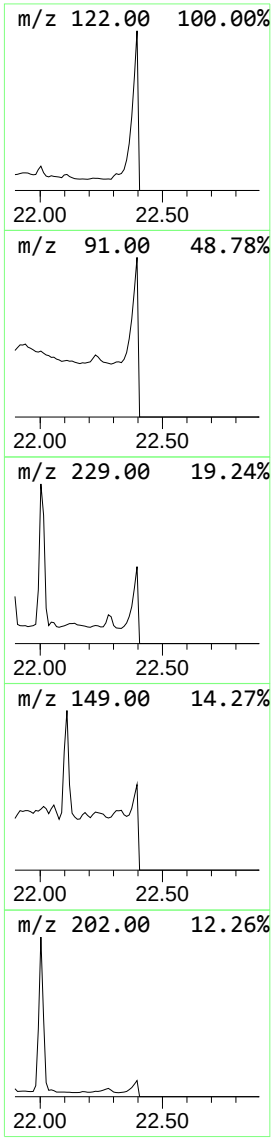
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2-Cyclopropyl-1H-indole, TMS Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.396	0.09 ng	51081	Perylene-d12	23.481

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclopropyl-1H-indole, TMS	229	C14H19NSi	1000494-09-0	9
2	1,3-Dioxolane, 2-(3-bromophenyl)-	228	C9H9BrO2	017789-14-9	7
3	1,3-Dioxolane, 2-(2-bromophenyl)-	228	C9H9BrO2	034824-58-3	7
4	[1,2,4]Triazolo[4,3-b]pyridazin-...	229	C11H11N5O	1000338-08-6	4
5	Benzene, 1-methoxy-3-methyl-	122	C8H10O	000100-84-5	4





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BNA_N
ClientSampleId :
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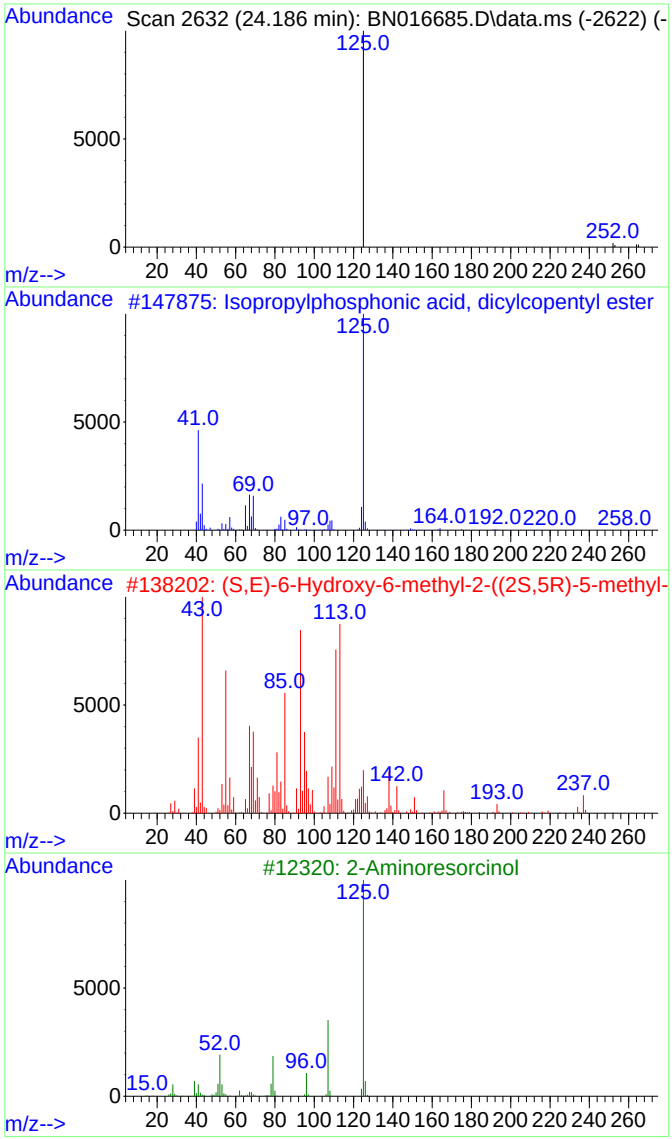
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

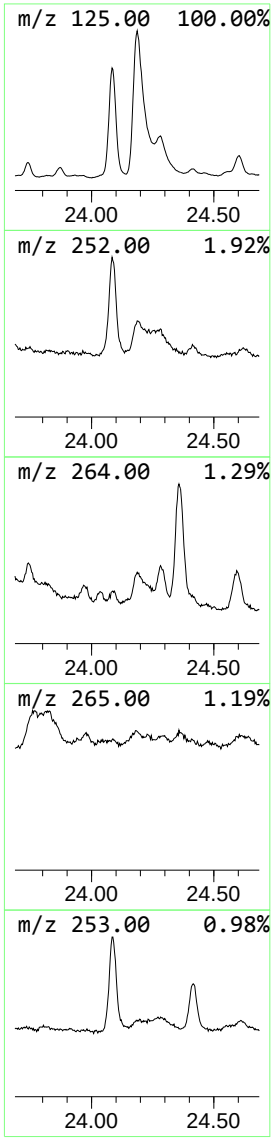
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Isopropylphosphonic acid, d... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.186	0.06 ng	31633	Perylene-d12	23.481

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropylphosphonic acid, dicylc...	260	C13H25O3P	1000273-18-4	3
2	(S,E)-6-Hydroxy-6-methyl-2-((2S,...	252	C15H24O3	088243-64-5	3
3	2-Aminoresorcinol	125	C6H7NO2	003163-15-3	2
4	2-Amino-4-hydroxy-6-methylpyrimi...	125	C5H7N3O	003977-29-5	2
5	2(1H)-Pyrimidinone, 4-amino-5-me...	125	C5H7N3O	000554-01-8	2





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
Data File : BN016685.D
Acq On : 02 Oct 2021 23:45
Operator : CG/JU
Sample : M3892-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
S-829-J-SO-3-3.5-092221

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Hydrogen isocya...	3.512	0.1	ng	24478	1	7.643	94584	0.4
beta-Alanine amide	3.825	0.0	ng	11062	1	7.643	94584	0.4
Butane	4.821	0.3	ng	70543	1	7.643	94584	0.4
Isobutane	5.042	0.0	ng	11561	1	7.643	94584	0.4
3-Hexanone, 2,5...	7.938	0.1	ng	17852	1	7.643	94584	0.4
Propane, 1-isot...	10.191	0.1	ng	57720	2	10.407	173562	0.4
1,3-Dioxolane, ...	20.474	0.1	ng	49085	5	21.228	291118	0.4
1,2,3-Benzenetr...	20.973	0.1	ng	60193	5	21.228	291118	0.4
Pyrazinamine, 3...	21.090	0.2	ng	133174	5	21.228	291118	0.4
acetamide, N-[5...	21.875	0.1	ng	106592	5	21.228	291118	0.4
5-bromo-2-hydro...	22.003	0.1	ng	50524	5	21.228	291118	0.4
2-Cyclopropyl-1...	22.396	0.1	ng	51081	6	23.481	229041	0.4
Isopropylphosph...	24.186	0.1	ng	31633	6	23.481	229041	0.4