LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_G\Methods\SFAM-EPA-BG120821.M
Title : SVOA CALIBRATION

Signal : TIC: BG051419.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.301	132	138	151	rVB	94217	158597	0.25%	0.202%
2	5.652	363	368	373	rVB	47844	76687	0.12%	0.098%
3	5.787	384	391	401	rBV	91786	205181	0.32%	0.261%
4	6.163	447	455	460	rVV2	53079	99566	0.16%	0.127%
5	6.845	564	571	579	rVB	103402	173914	0.27%	0.222%
6	7.509	678	684	690	rBV	100323	191321	0.30%	0.244%
7	7.579	691	696	708	rVB	41221	92849	0.15%	0.118%
8	8.190	793	800	805	rBV	96991	172598	0.27%	0.220%
9	9.095	947	954	966	rBV	283362	648881	1.02%	0.827%
10	10.035	1108	1114	1128	rVB	168956	320591	0.51%	0.409%
11	10.194	1133	1141	1158	rBV3	28757	100733	0.16%	0.128%
12	10.358	1158	1169	1174	rBV2	39841	88701	0.14%	0.113%
13	10.529	1188	1198	1208	rVV4	22916	68691	0.11%	0.088%
14	11.016	1271	1281	1285	rBV	151667	286606	0.45%	0.365%
15	11.063	1285	1289	1296	rVB	102917	188271	0.30%	0.240%
16	11.328	1296	1334	1338	rBV	10166274	63436645	100.00%	80.842%
17	12.485	1524	1531	1534	rBV	76068	138399	0.22%	0.176%
18	12.550	1534	1542	1555	rVV	3386697	5883679	9.27%	7.498%
19	13.278	1659	1666	1673	rBV	82342	141468	0.22%	0.180%
20	14.054	1793	1798	1804	rVB	69894	106161	0.17%	0.135%
21	14.183	1815	1820	1823	rBV	41602	74043	0.12%	0.094%
22	14.688	1900	1906	1912	rVB	64354	101392	0.16%	0.129%
23	14.824	1918	1929	1937	rVB	227304	373570	0.59%	0.476%
24	14.970	1947	1954	1965	rBV	62916	99437	0.16%	0.127%
25	15.570	2049	2056	2060	rBV2	83589	128808	0.20%	0.164%
26	15.646	2060	2069	2079	rVV	1575168	3374476	5.32%	4.300%
27	15.811	2091	2097	2102	rBV2	71736	114538	0.18%	0.146%
28	17.573	2390	2397	2408	rVB	287753	447692	0.71%	0.571%
29	17.673	2408	2414	2419	rBV	49797	77482	0.12%	0.099%
30	19.953	2794	2802	2807	rBV3	59123	103250	0.16%	0.132%
31	20.834	2948	2952	2960	rBV	54761	76714	0.12%	0.098%
32	21.874	3122	3129	3138	rVB	262939	467761	0.74%	0.596%
33	25.276	3698	3708	3723	rVB3	144505	451486	0.71%	0.575%

Sum of corrected areas: 78470188

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
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Acq On : 8 Dec 2021 23:09
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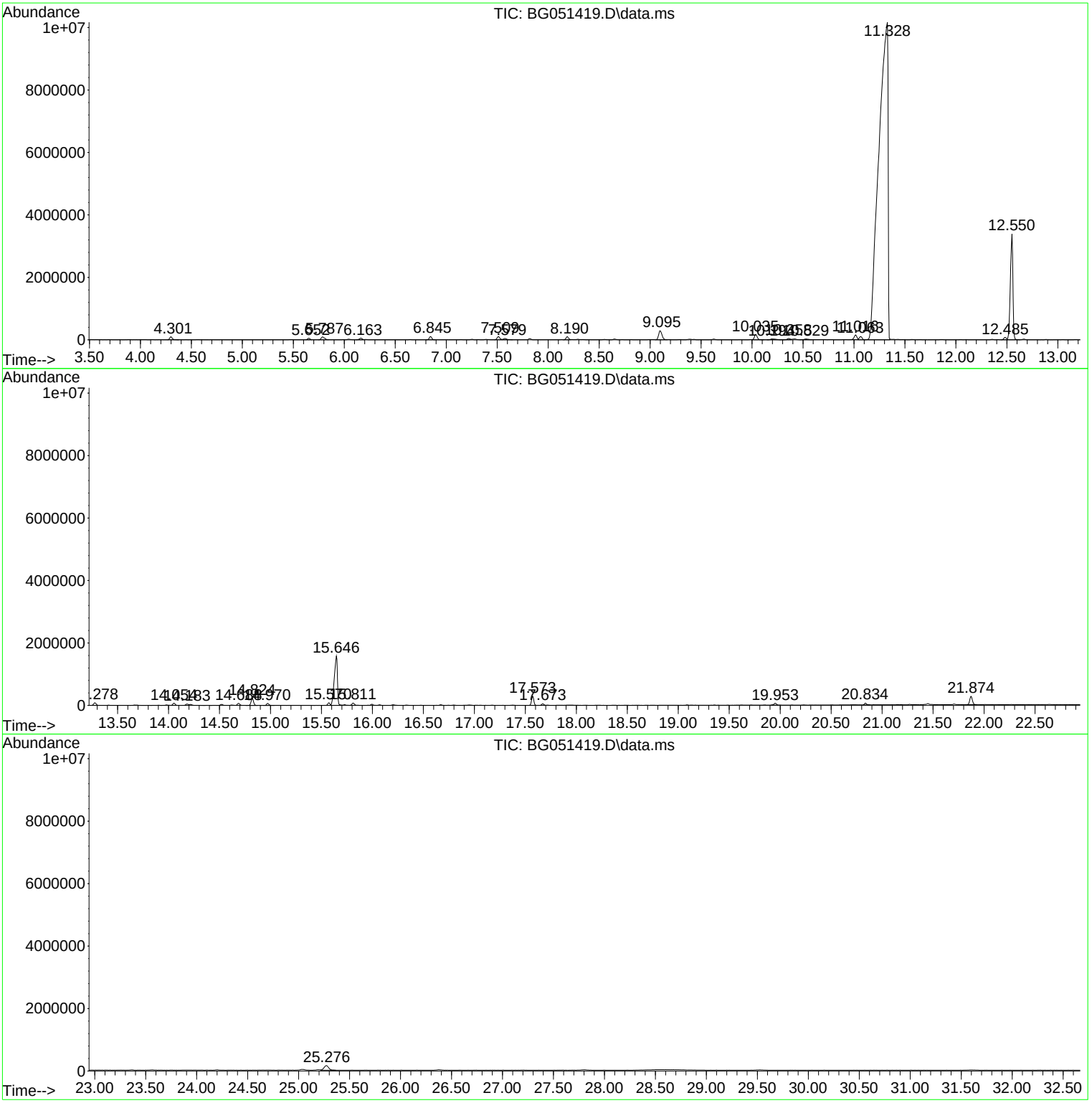
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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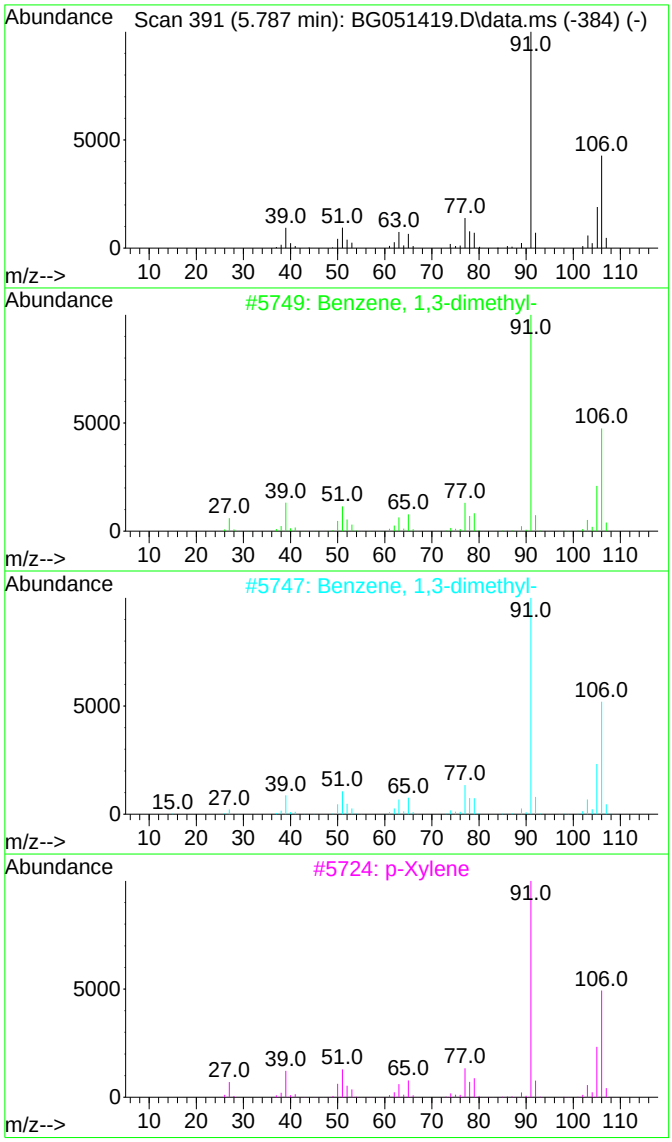
Quant Title : SVOA CALIBRATION

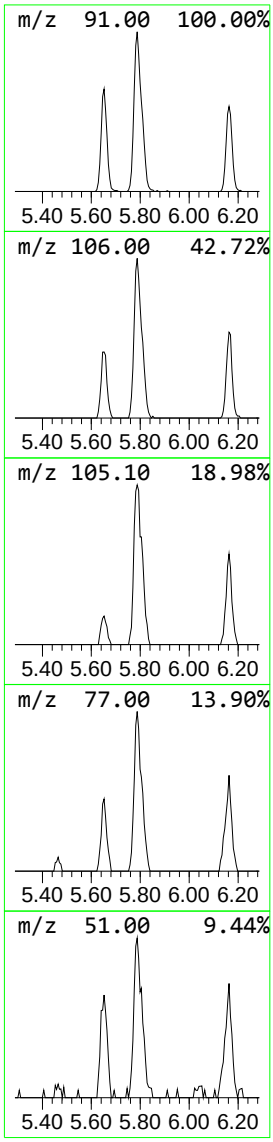
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.787	23.78 ng/ul	205181	1,4-Dichlorobenzene-d4	8.190

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	o-Xylene	106	C8H10	000095-47-6	95
5	o-Xylene	106	C8H10	000095-47-6	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

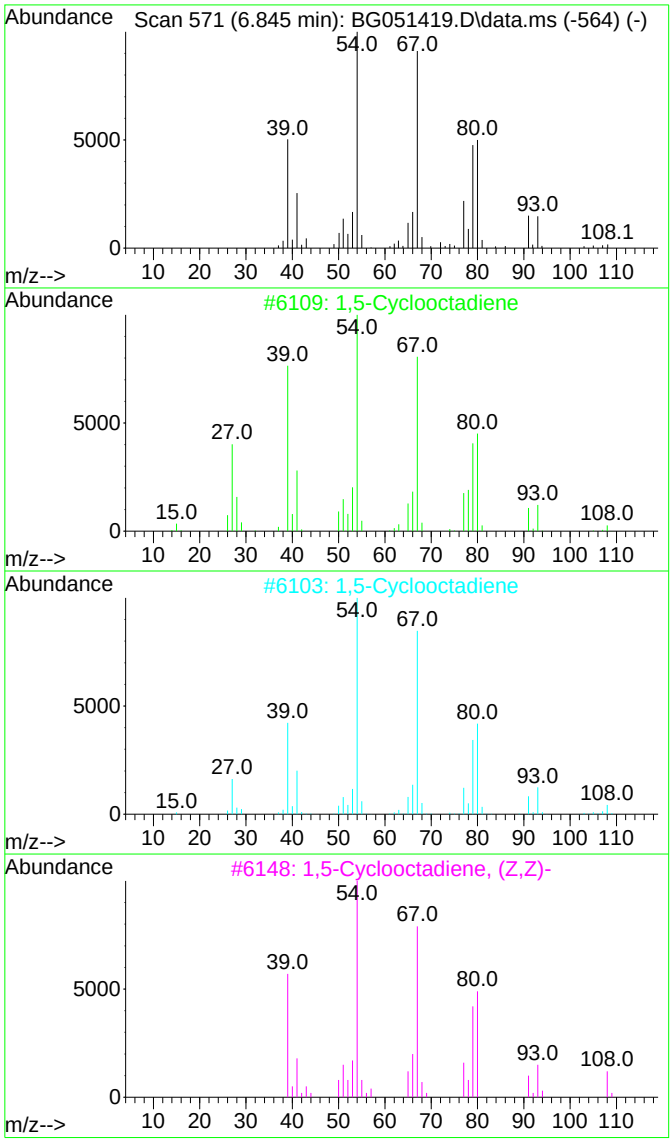
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

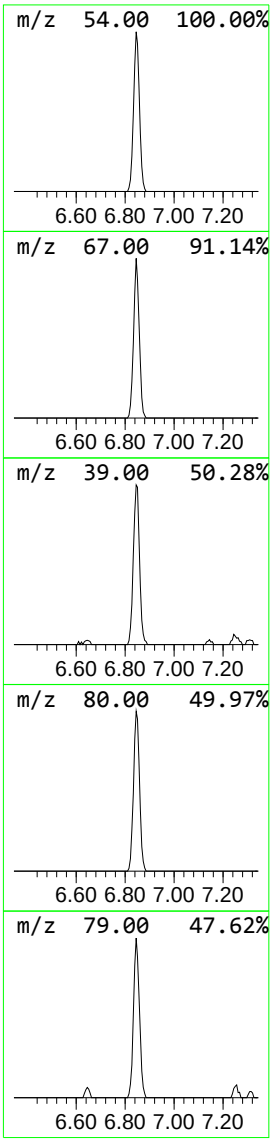
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,5-Cyclooctadiene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.845	20.15 ng/ul	173914	1,4-Dichlorobenzene-d4	8.190

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Cyclooctadiene	108	C8H12	000111-78-4	94
2	1,5-Cyclooctadiene	108	C8H12	000111-78-4	90
3	1,5-Cyclooctadiene, (Z,Z)-	108	C8H12	001552-12-1	90
4	1,5-Cyclooctadiene	108	C8H12	000111-78-4	90
5	1,5-Cyclooctadiene, (E,Z)-	108	C8H12	005259-71-2	86





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Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

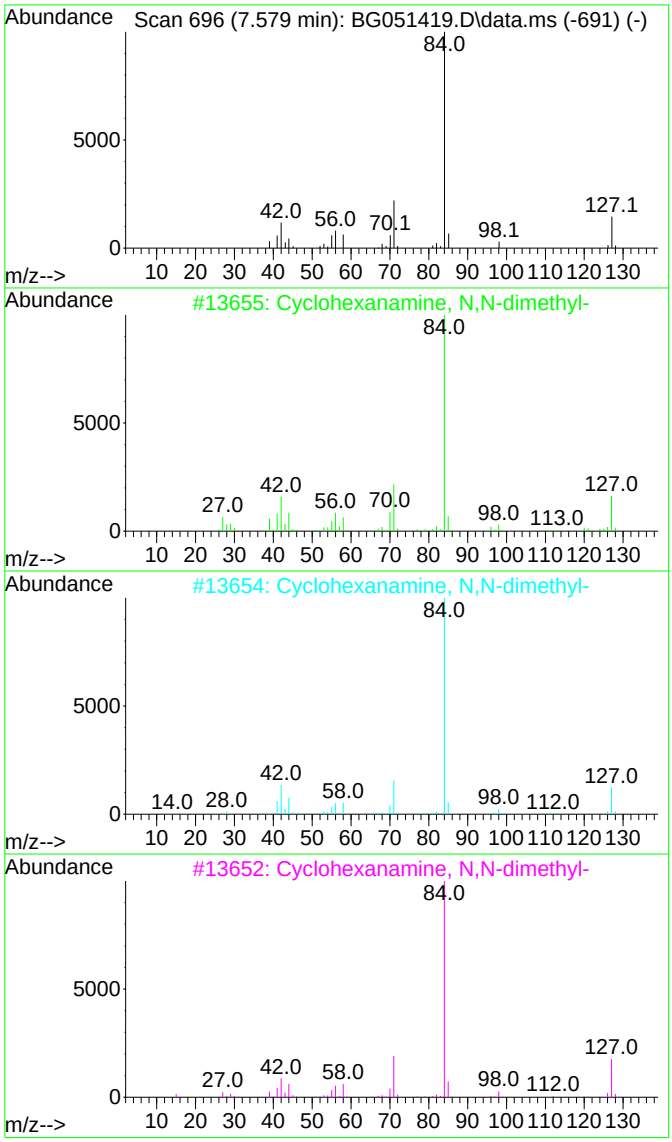
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

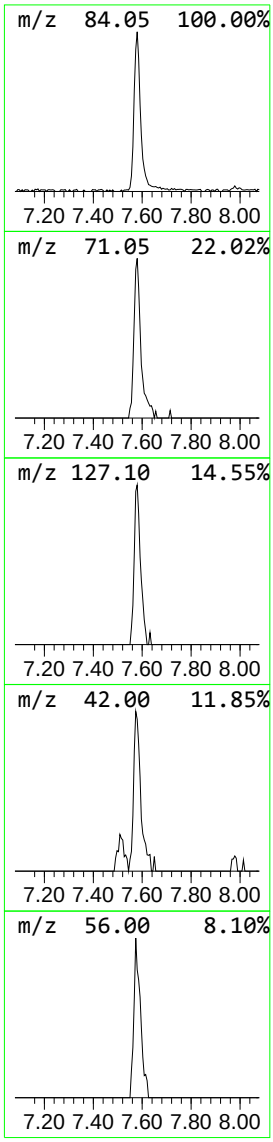
****

Peak Number 6 Cyclohexanamine, N,N-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.579	10.76 ng/ul	92849	1,4-Dichlorobenzene-d4	8.190

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N,N-dimethyl-	127	C8H17N	000098-94-2	91
2	Cyclohexanamine, N,N-dimethyl-	127	C8H17N	000098-94-2	90
3	Cyclohexanamine, N,N-dimethyl-	127	C8H17N	000098-94-2	90
4	3-Amino-1,2,4-triazole-5-carboxy...	128	C3H4N4O2	003641-13-2	50
5	1H-Pyrrole-1-ethanamine, tetrahy...	128	C7H16N2	1000351-26-6	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

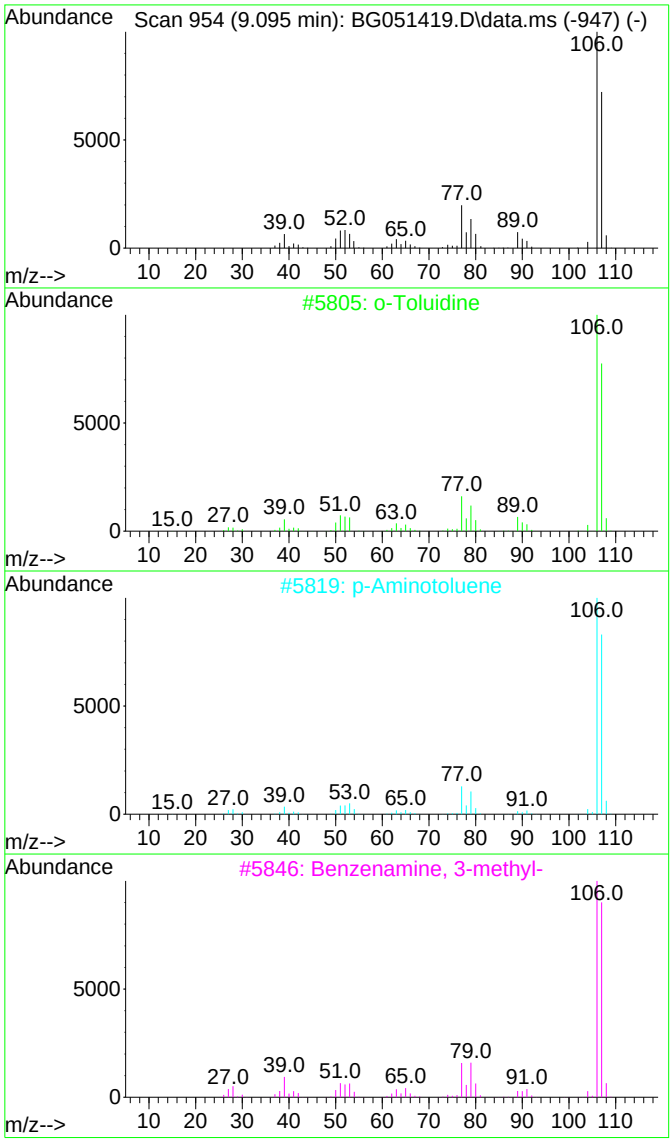
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

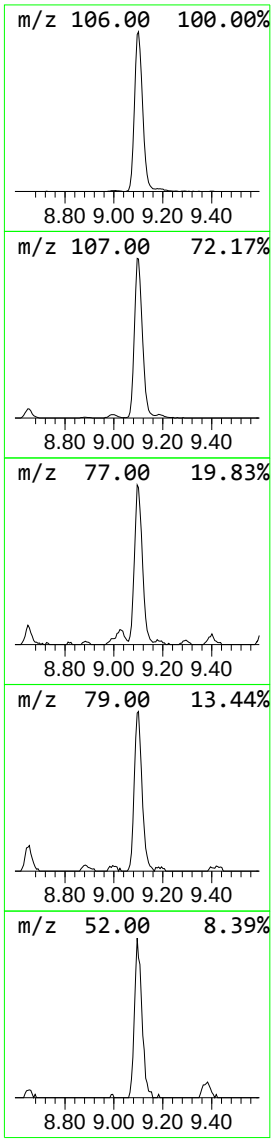
****

Peak Number 7 o-Toluidine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.095	75.19 ng/ul	648881	1,4-Dichlorobenzene-d4	8.190

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Toluidine	107	C7H9N	000095-53-4	97
2	p-Aminotoluene	107	C7H9N	000106-49-0	94
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	93
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
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Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

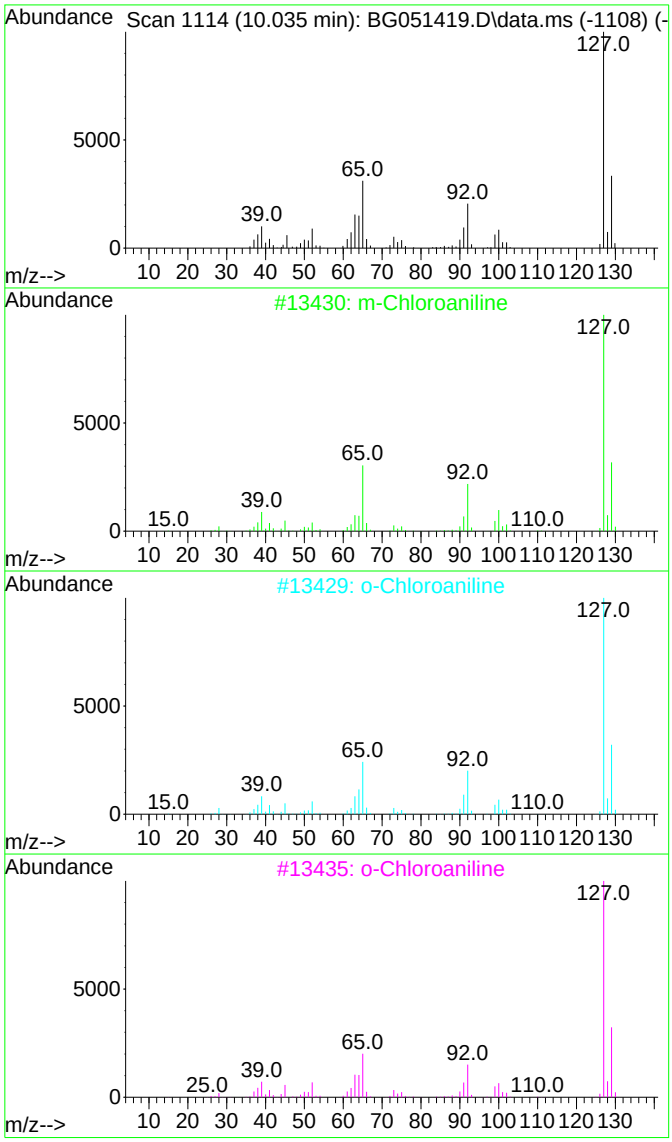
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

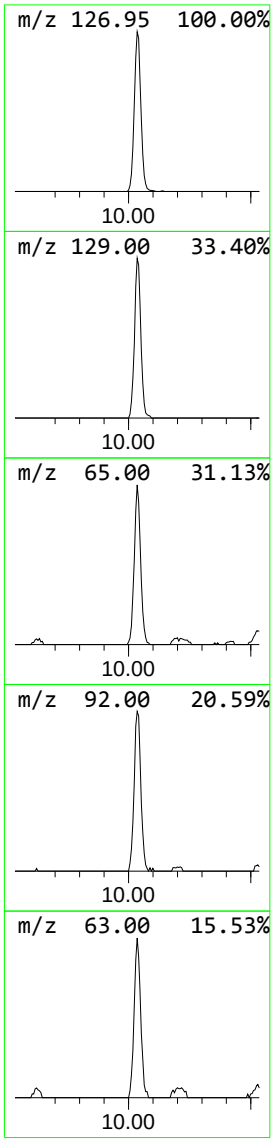
****

Peak Number 8 m-Chloroaniline Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.035	22.37 ng/ul	320591	Naphthalene-d8	11.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Chloroaniline	127	C6H6ClN	000108-42-9	97
2	o-Chloroaniline	127	C6H6ClN	000095-51-2	96
3	o-Chloroaniline	127	C6H6ClN	000095-51-2	95
4	p-Chloroaniline	127	C6H6ClN	000106-47-8	95
5	m-Chloroaniline	127	C6H6ClN	000108-42-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

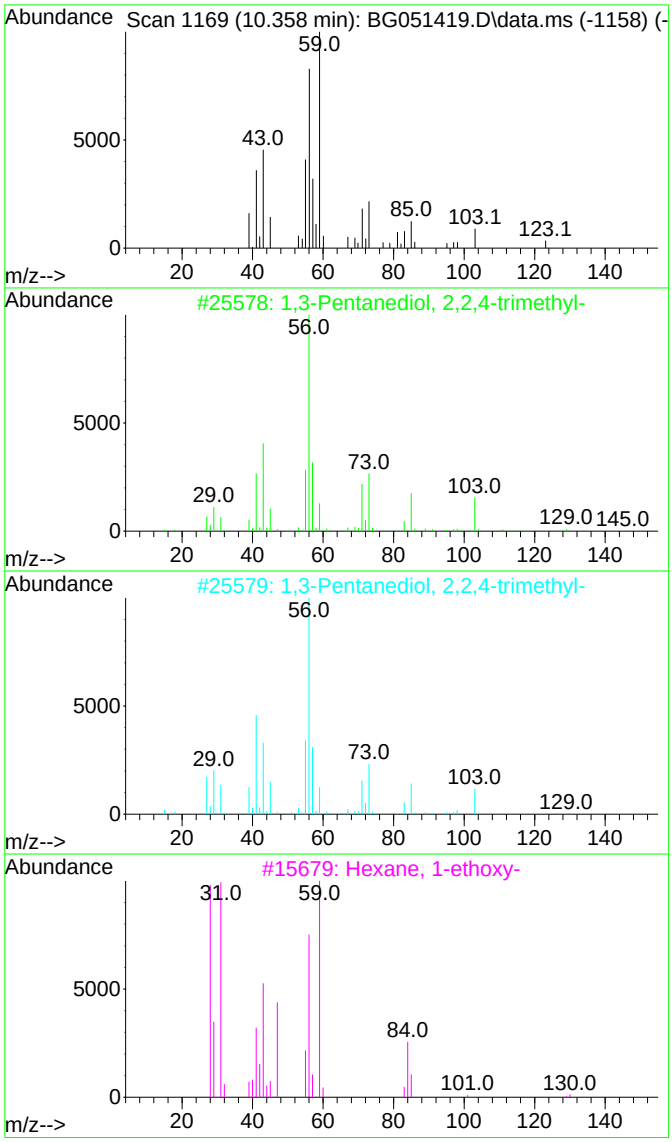
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

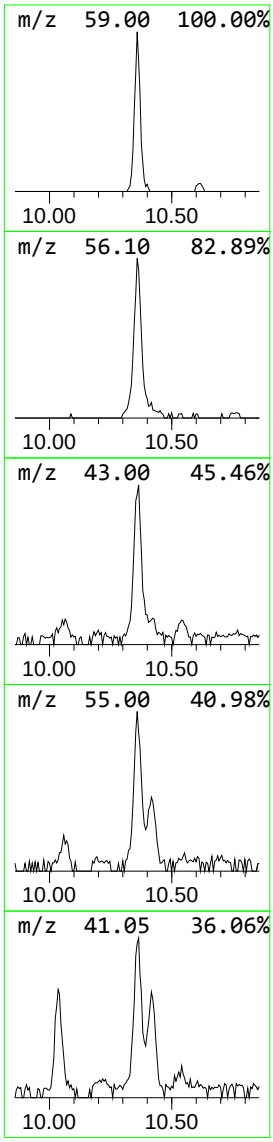
****

Peak Number 9 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.358	6.19 ng/ul	88701	Naphthalene-d8	11.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	47
2	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	47
3	Hexane, 1-ethoxy-	130	C8H18O	005756-43-4	45
4	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	43
5	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	43





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Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

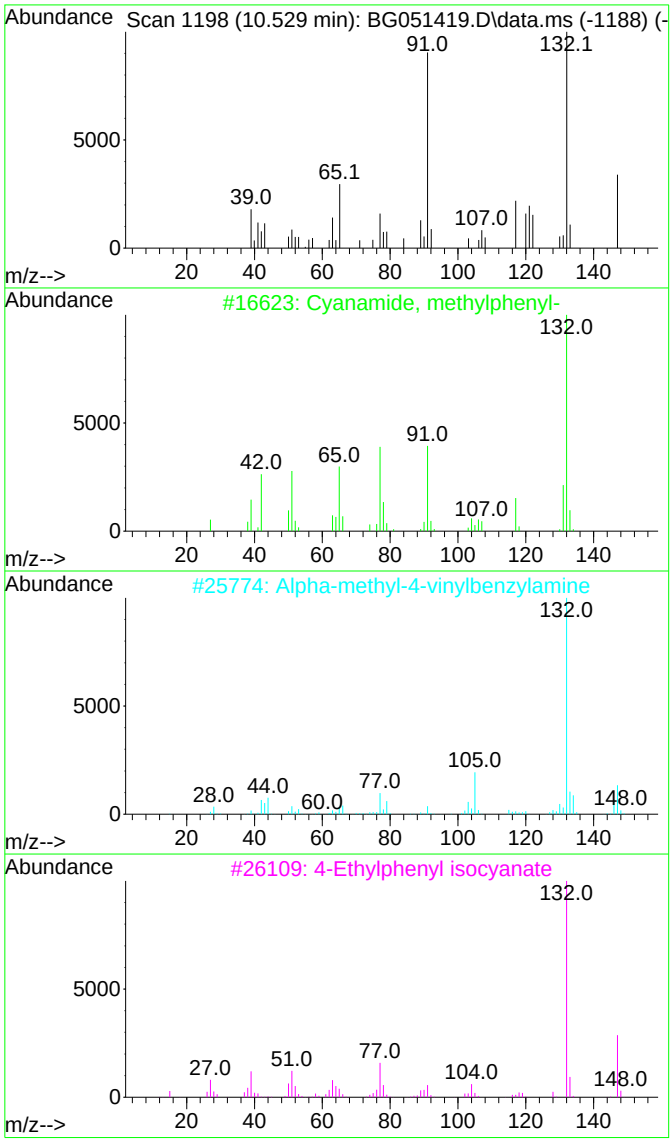
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

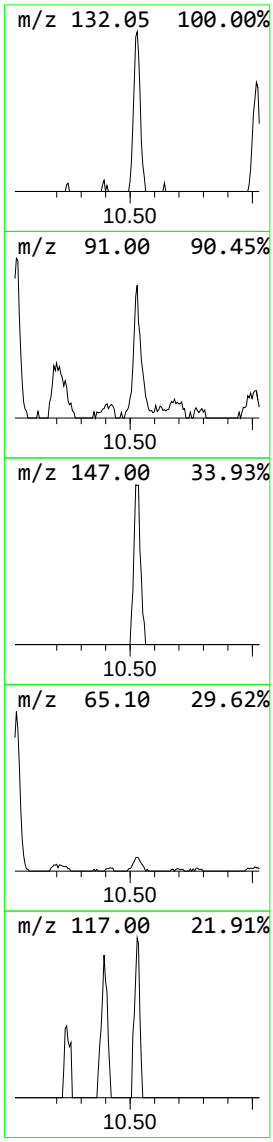
****

Peak Number 10 Cyanamide, methylphenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.529	4.79 ng/ul	68691	Naphthalene-d8	11.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyanamide, methylphenyl-	132	C8H8N2	018773-77-8	50
2	Alpha-methyl-4-vinylbenzylamine	147	C10H13N	019303-08-3	46
3	4-Ethylphenyl isocyanate	147	C9H9NO	023138-50-3	43
4	Methyl-3-phenylaziridine, 2-	133	C9H11N	007763-71-5	43
5	1-Aminoindan	133	C9H11N	034698-41-4	43





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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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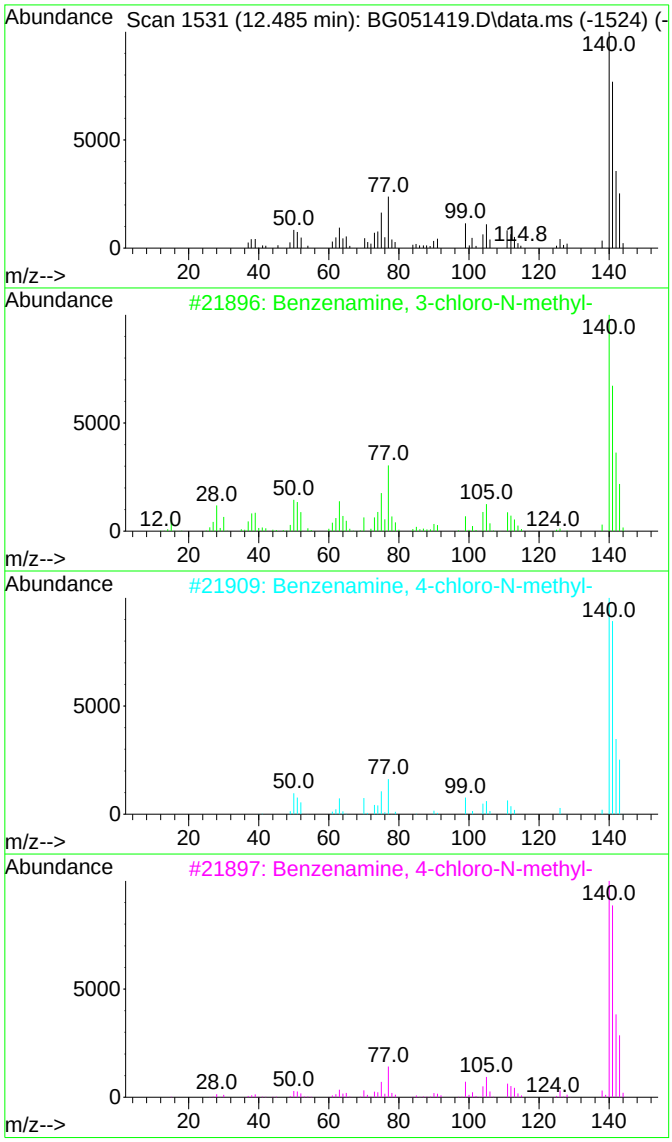
Quant Title : SVOA CALIBRATION

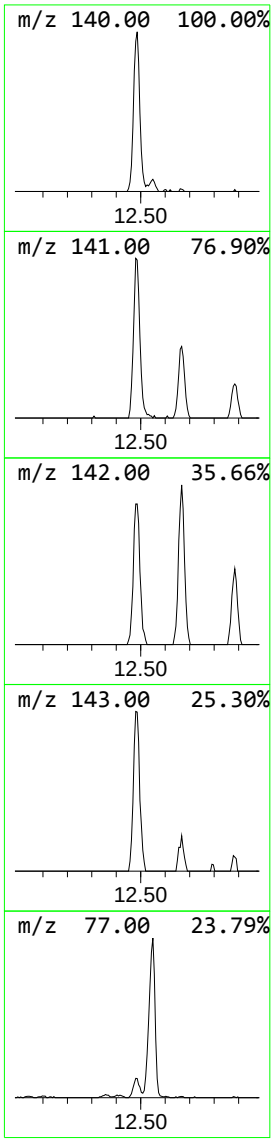
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzenamine, 3-chloro-N-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.485	9.66 ng/ul	138399	Naphthalene-d8	11.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-N-methyl-	141	C7H8ClN	007006-52-2	97
2	Benzenamine, 4-chloro-N-methyl-	141	C7H8ClN	000932-96-7	95
3	Benzenamine, 4-chloro-N-methyl-	141	C7H8ClN	000932-96-7	95
4	Benzenamine, 3-chloro-N-methyl-	141	C7H8ClN	007006-52-2	93
5	Benzenamine, 2-chloro-N-methyl-	141	C7H8ClN	000932-32-1	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
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Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

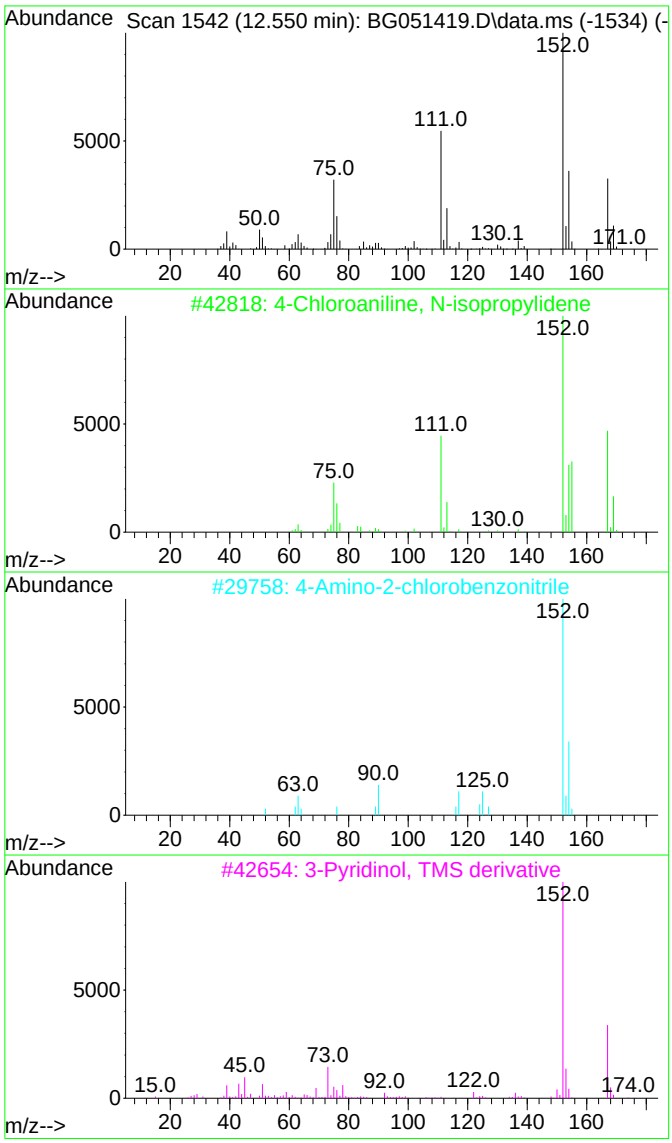
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

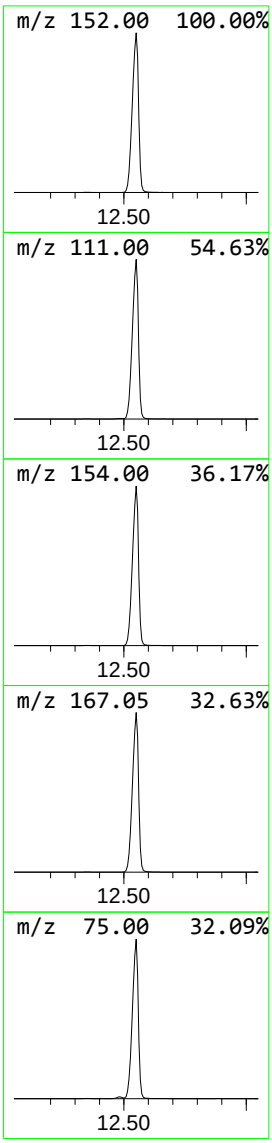
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 4-Chloroaniline, N-isopropy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.550	410.58 ng/ul	5883680	Naphthalene-d8	11.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloroaniline, N-isopropylidene	167	C9H10ClN	040938-43-0	91
2	4-Amino-2-chlorobenzonitrile	152	C7H5ClN2	020925-27-3	43
3	3-Pyridinol, TMS derivative	167	C8H13NOSi	041571-88-4	38
4	5-Aminopyrimidine, N-trimethylsi...	167	C7H13N3Si	1000493-88-4	38
5	2-Chlorobenzimidazole	152	C7H5ClN2	004857-06-1	38





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Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

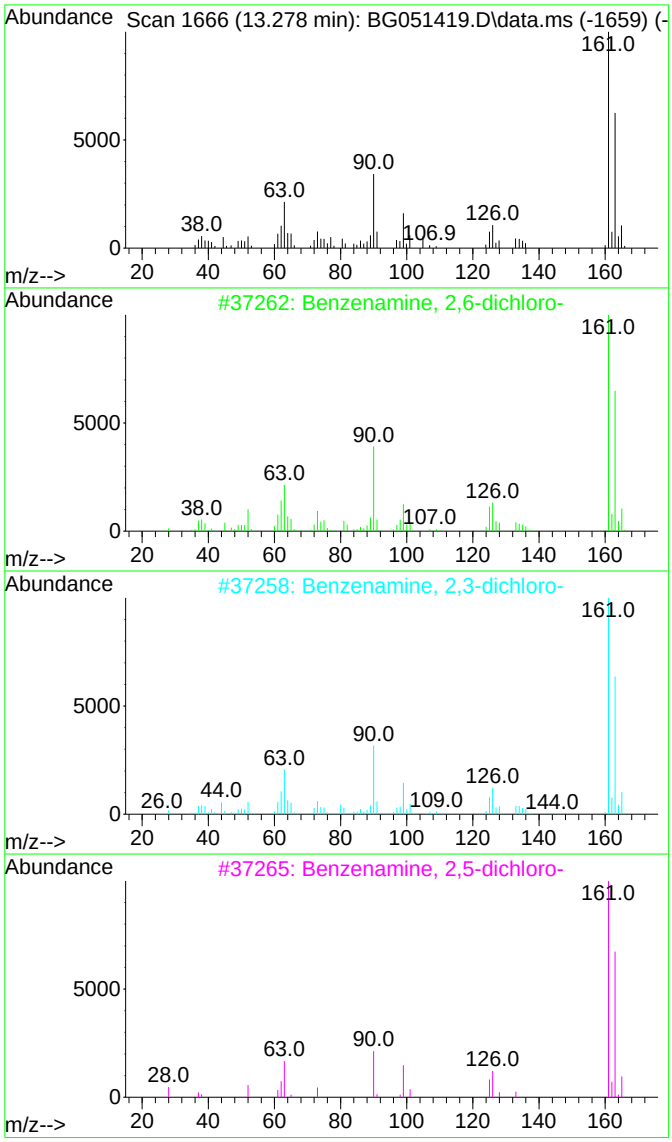
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

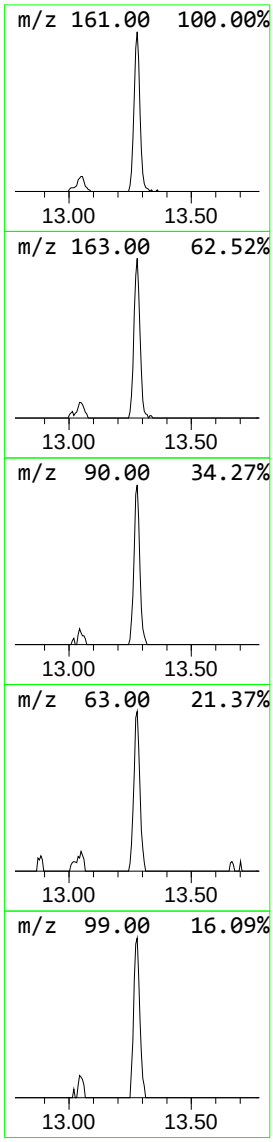
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzenamine, 2,6-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.278	7.57 ng/ul	141468	Acenaphthene-d10	14.824

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	98
2	Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
3	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
4	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	95
5	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

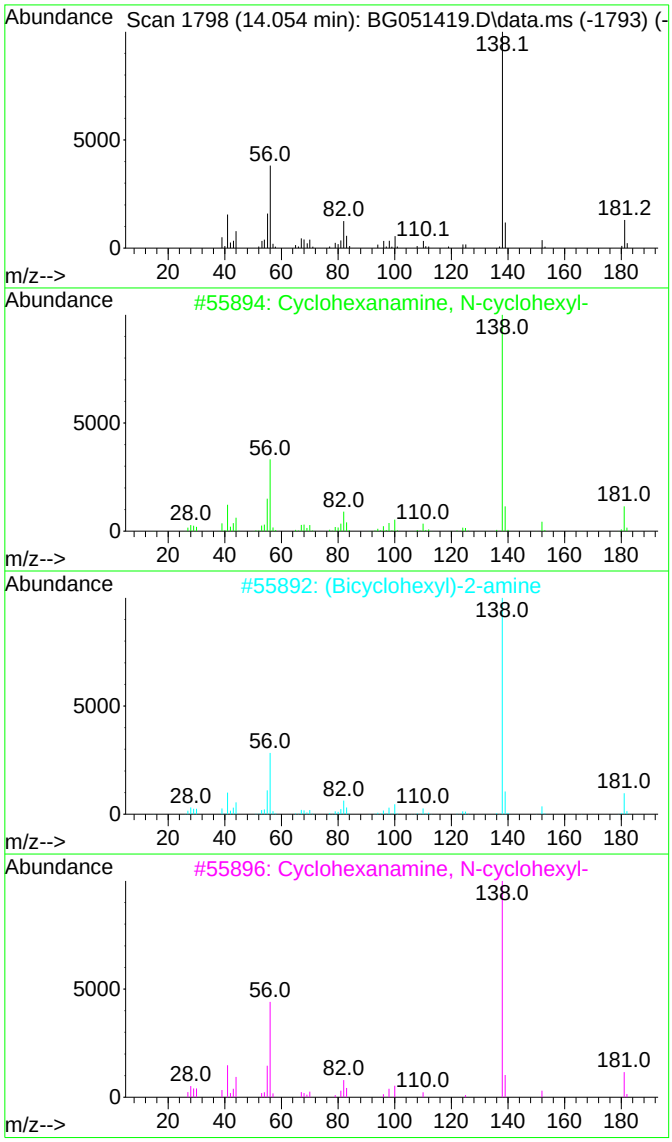
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

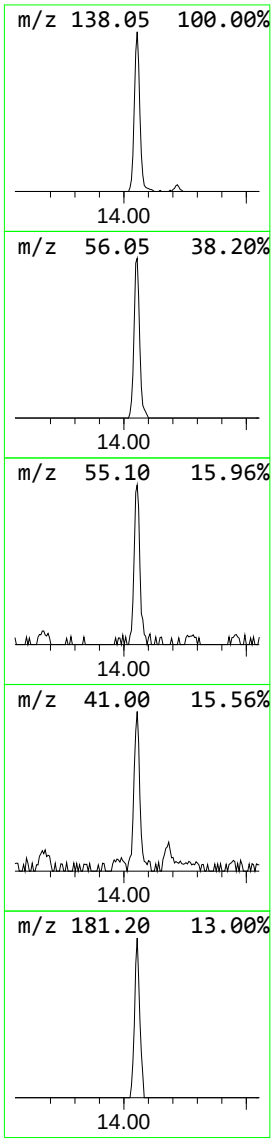
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Cyclohexanamine, N-cyclohexyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.054	5.68 ng/ul	106161	Acenaphthene-d10	14.824

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
2	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	96
3	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	95
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

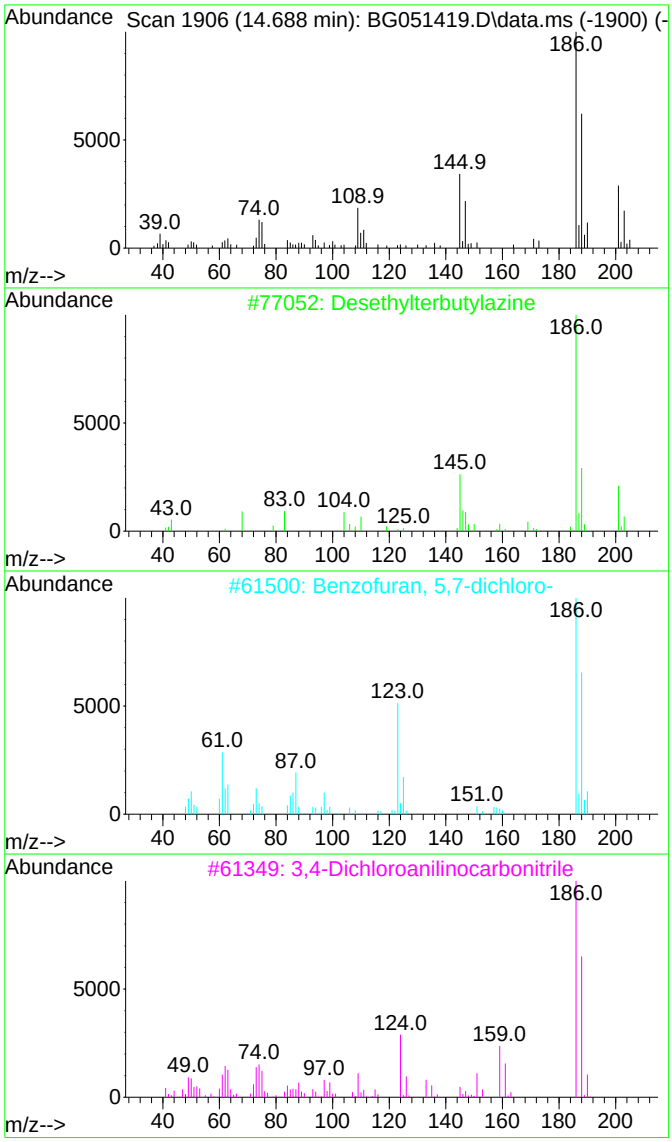
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

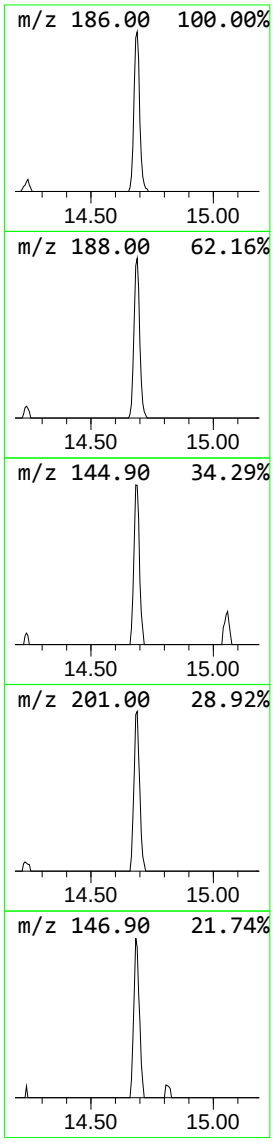
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Desethylterbutylazine Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.688	5.43 ng/ul	101392	Acenaphthene-d10	14.824

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Desethylterbutylazine	201	C7H12ClN5	030125-63-4	53
2	Benzofuran, 5,7-dichloro-	186	C8H4Cl2O	023145-06-4	47
3	3,4-Dichloroanilinocarbonitrile	186	C7H4Cl2N2	018995-48-7	43
4	(3-Chloro-5-fluorophenyl)diethyl...	201	C10H13ClFN	1000306-63-3	40
5	2-Cyanoamino-4-hydroxyquinazoline	186	C9H6N4O	112656-43-6	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

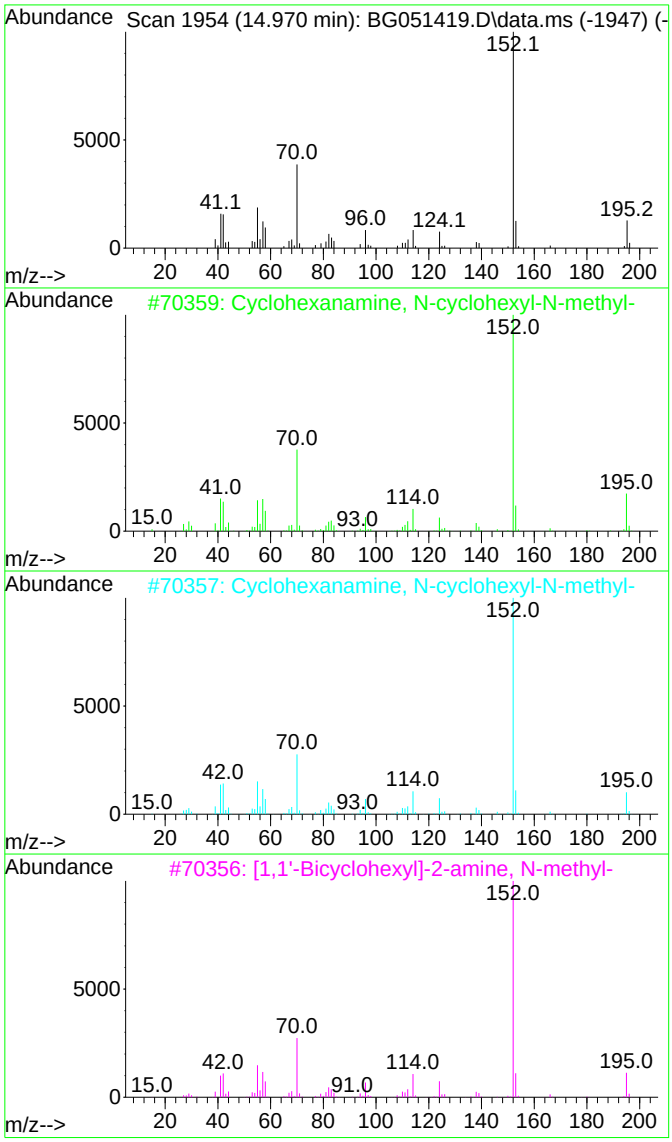
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

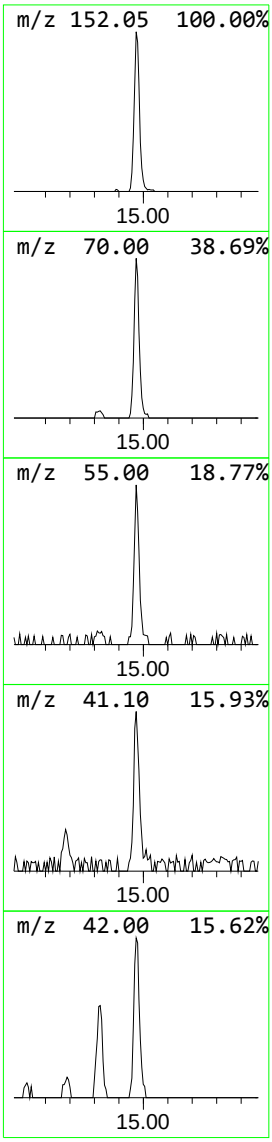
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Cyclohexanamine, N-cyclohex... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.970	5.32 ng/ul	99437	Acenaphthene-d10	14.824

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-N-...	195	C13H25N	007560-83-0	96
2	Cyclohexanamine, N-cyclohexyl-N-...	195	C13H25N	007560-83-0	96
3	[1,1'-Bicyclohexyl]-2-amine, N-m...	195	C13H25N	1010452-63-0	96
4	Cyclohexanamine, N-cyclohexyl-N-...	195	C13H25N	007560-83-0	94
5	Thieno[3,2-d]pyrimidin-4(3H)-one	152	C6H4N2OS	016234-10-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

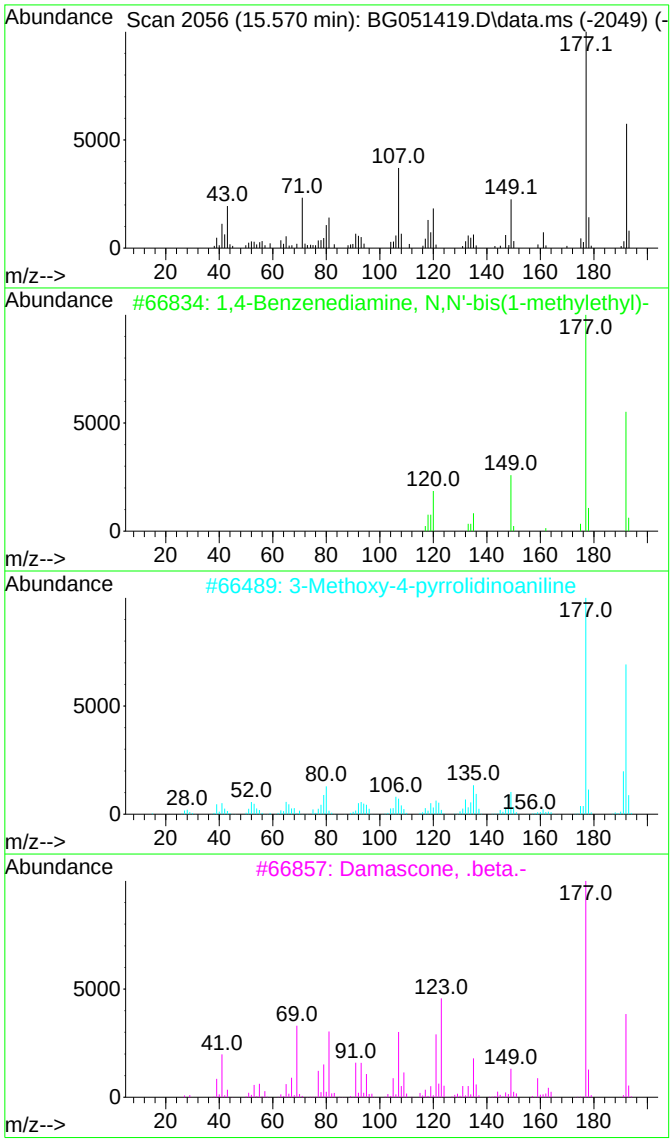
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA CALIBRATION

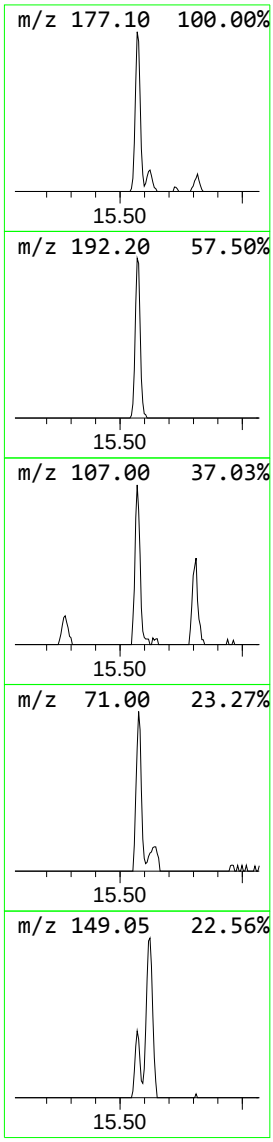
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1,4-Benzenediamine, N,N'-bi... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.570	6.90 ng/ul	128808	Acenaphthene-d10	14.824

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenediamine, N,N'-bis(1-m...	192	C12H20N2	004251-01-8	94
2	3-Methoxy-4-pyrrolidinoaniline	192	C11H16N2O	016089-42-2	58
3	Damascone, .beta.-	192	C13H20O	023726-91-2	58
4	5-(4,5-Dihydro-3H-pyrrol-2-ylmet...	192	C11H16N2O	051430-87-6	53
5	N-[4-(Ethyl-methyl-amino)-phenyl...	192	C11H16N2O	099981-45-0	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
Data File : BG051419.D
Acq On : 8 Dec 2021 23:09
Operator : CG/JU
Sample : M4960-11 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKS4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
Quant Title : SVOA 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
Benzene, 1,3-di...	5.787	23.8	ng/ul	205181	1	8.190	172598	20.0
1,5-Cyclooctadiene	6.845	20.1	ng/ul	173914	1	8.190	172598	20.0
Cyclohexanamine...	7.579	10.8	ng/ul	92849	1	8.190	172598	20.0
o-Toluidine	9.095	75.2	ng/ul	648881	1	8.190	172598	20.0
m-Chloroaniline	10.035	22.4	ng/ul	320591	2	11.016	286606	20.0
unknown-01	10.358	6.2	ng/ul	88701	2	11.016	286606	20.0
Cyanamide, meth...	10.529	4.8	ng/ul	68691	2	11.016	286606	20.0
Benzenamine, 3-...	12.485	9.7	ng/ul	138399	2	11.016	286606	20.0
4-Chloroaniline...	12.550	410.6	ng/ul	5883680	2	11.016	286606	20.0
Benzenamine, 2,...	13.278	7.6	ng/ul	141468	3	14.824	373570	20.0
Cyclohexanamine...	14.054	5.7	ng/ul	106161	3	14.824	373570	20.0
Desethylterbuty...	14.688	5.4	ng/ul	101392	3	14.824	373570	20.0
Cyclohexanamine...	14.970	5.3	ng/ul	99437	3	14.824	373570	20.0
1,4-Benzenediam...	15.570	6.9	ng/ul	128808	3	14.824	373570	20.0