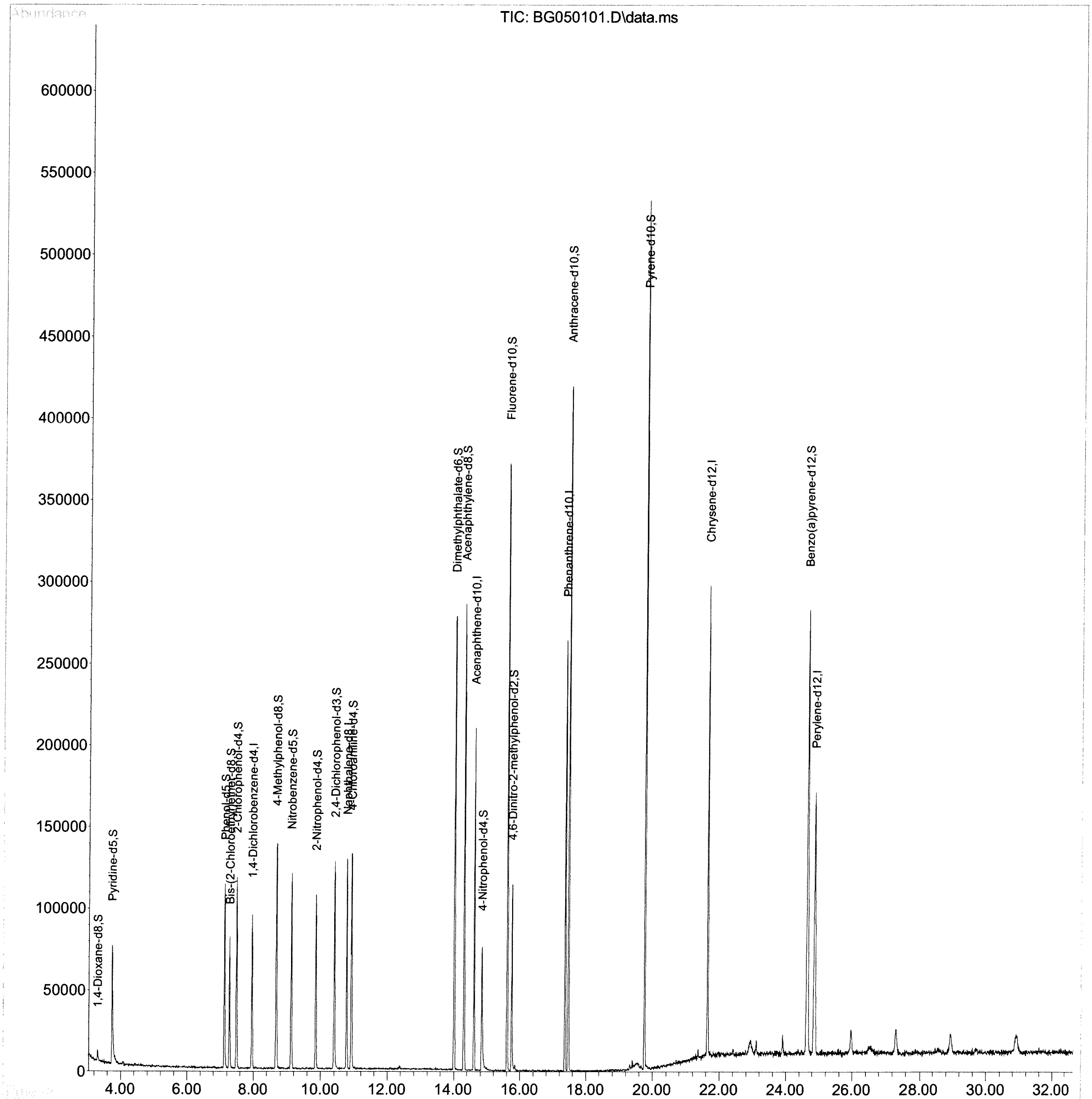
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050101.D
Acq On : 2 Sep 2021 17:47
Operator : CG/JU
Sample : PB138904BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Sep 02 18:33:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SBLK904

Manual Integrations APPROVED
mohammad
9/3/2021 3:31:42 PM



Quantitation Report (Qedit)

```
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050101.D
Acq On    :  2 Sep 2021  17:47
Operator  : CG/JU
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Instrument :
BNA_G
ClientSampleId :
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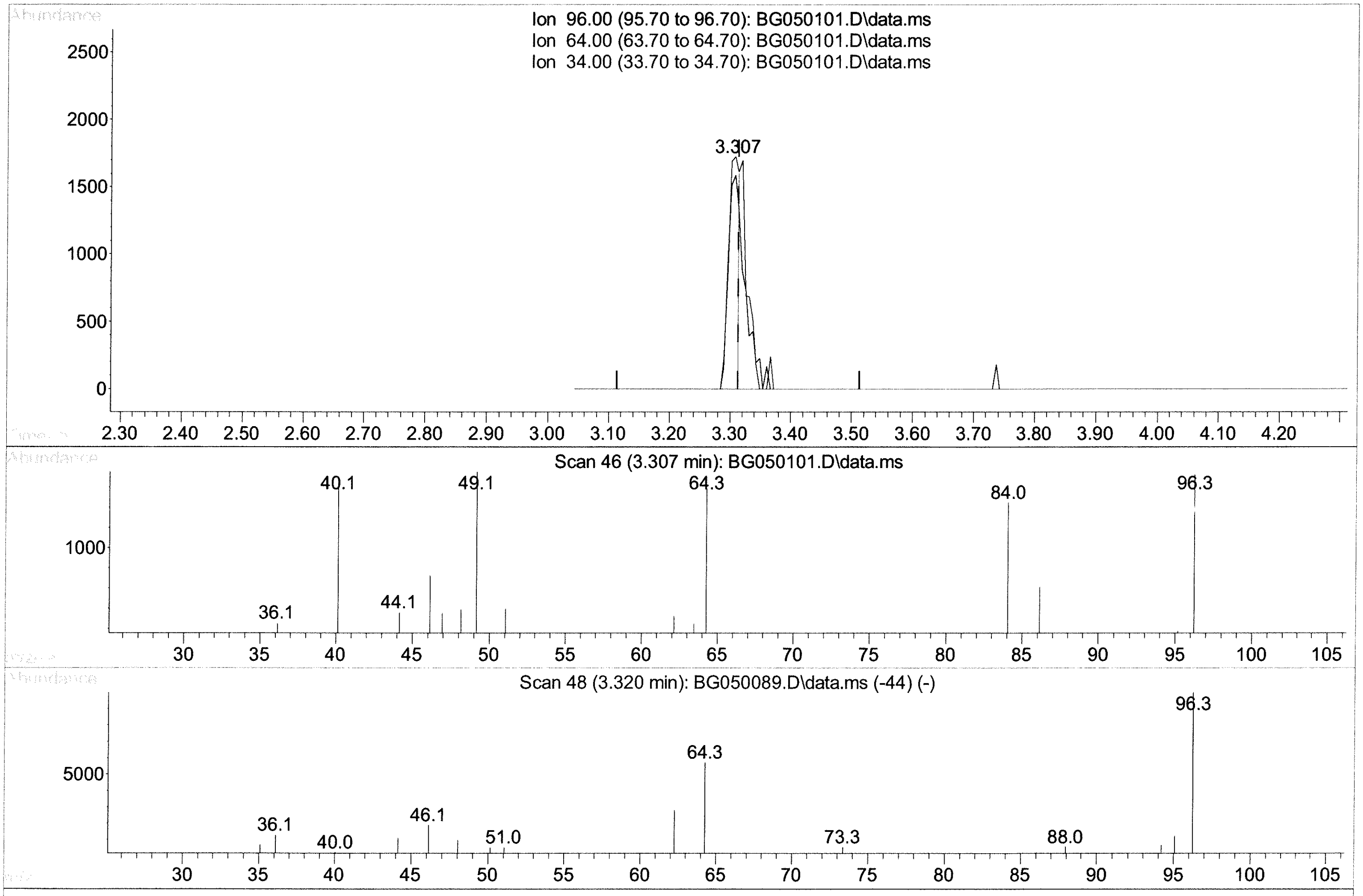
Manual Integrations APPROVED
mohammad
9/3/2021 3:31:42 PM



TIC: BG050101.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.307min (-0.006) 4.16 ng/uL

response 2160

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	91.99#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
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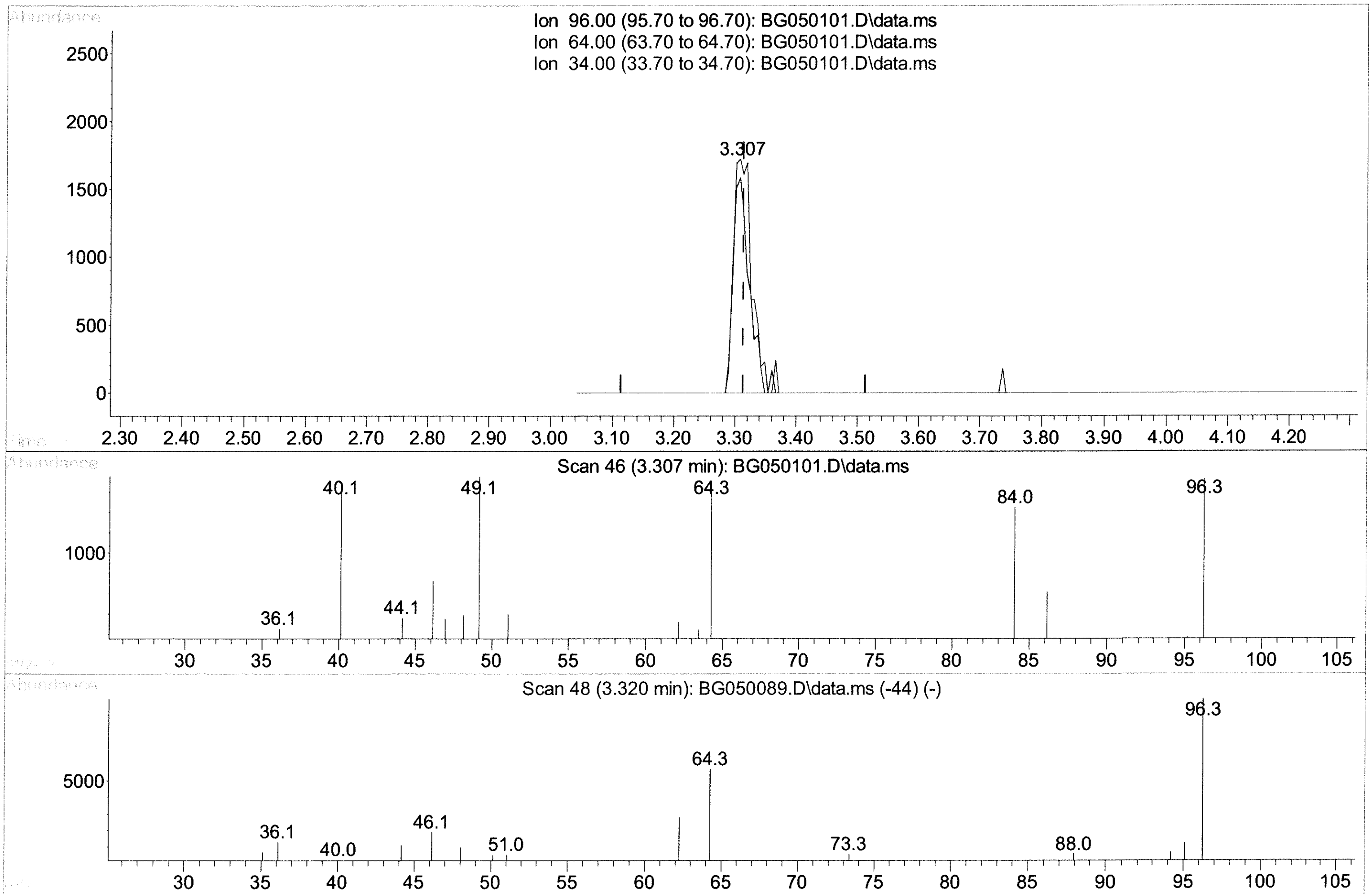
Manual Integrations APPROVED
mohammad
9/3/2021 3:31:42 PM

Quant Time: Sep 02 18:33:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration



TIC: BG050101.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.307min (-0.006) 6.72 ng/uL m JU 9/8/21

response 3485

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	91.99#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Data File : BG050101.D
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ALS Vial : 3 Sample Multiplier: 1

Quant Time: Sep 02 18:33:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration

Instrument : BNA_G
ClientSampleId : SBLK904

Manual Integrations APPROVED
mohammad
9/3/2021 3:31:42 PM

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	26758	20.000	ng/ul	0.00
20) Naphthalene-d8	10.774	136	108377	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	71601	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.371	188	162749	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.641	240	171905	20.000	ng/ul	0.00
88) Perylene-d12	24.866	264	173509	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	3485m	6.716	ng/uL	0.00 JY 9/8/21
4) Pyridine-d5	3.736	84	47724	30.534	ng/ul	0.00
7) Phenol-d5	7.126	99	68038	29.944	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.279	67	39843	31.563	ng/ul	0.00
11) 2-Chlorophenol-d4	7.484	132	52370	30.803	ng/ul	0.00
15) 4-Methylphenol-d8	8.671	113	53547	29.924	ng/ul	-0.01
21) Nitrobenzene-d5	9.123	128	27373	32.578	ng/ul	0.00
24) 2-Nitrophenol-d4	9.851	143	30539	30.524	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	49714	27.859	ng/ul	0.00
31) 4-Chloroaniline-d4	10.909	131	72989	30.051	ng/ul	-0.01
46) Dimethylphthalate-d6	14.016	166	187244	34.170	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	216314	33.668	ng/ul	0.00
54) 4-Nitrophenol-d4	14.851	143	21353	26.933	ng/ul	0.00
60) Fluorene-d10	15.608	176	151669	32.645	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.749	200	27556	26.318	ng/ul	0.00
73) Anthracene-d10	17.471	188	243934	33.378	ng/ul	0.00
81) Pyrene-d10	19.762	212	297469	32.022	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.643	264	290391	31.547	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed