

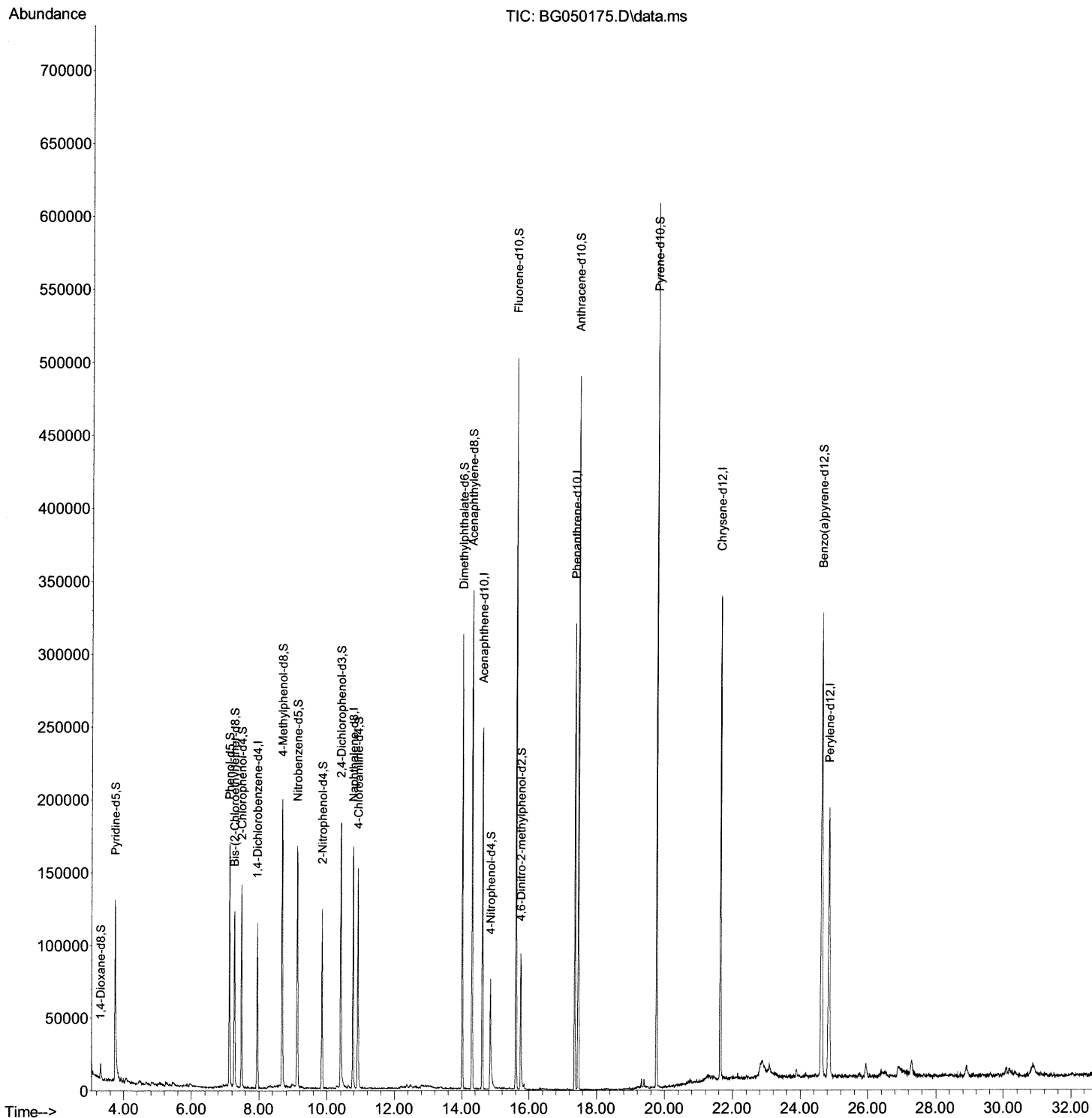
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050175.D
Acq On : 7 Sep 2021 21:21
Operator : CG/JU
Sample : PB138822BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK822

Manual Integrations
APPROVED

mohammad
9/9/2021 8:39:40 AM

Quant Time: Sep 08 01:49:17 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

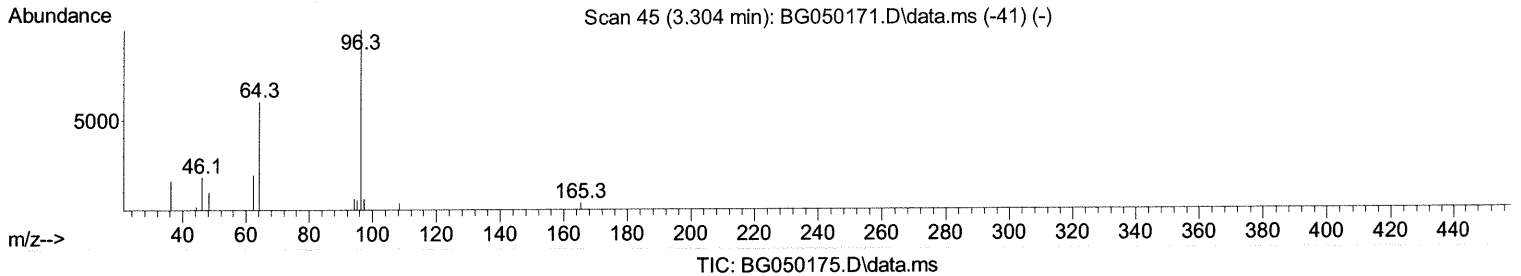
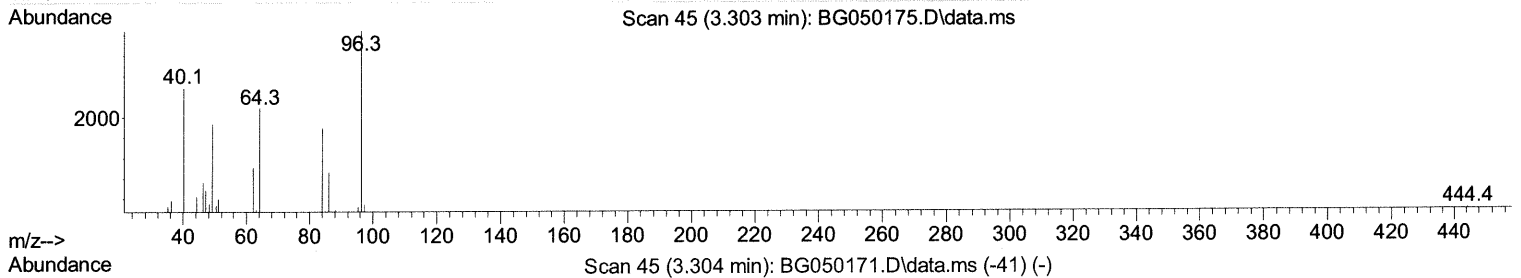
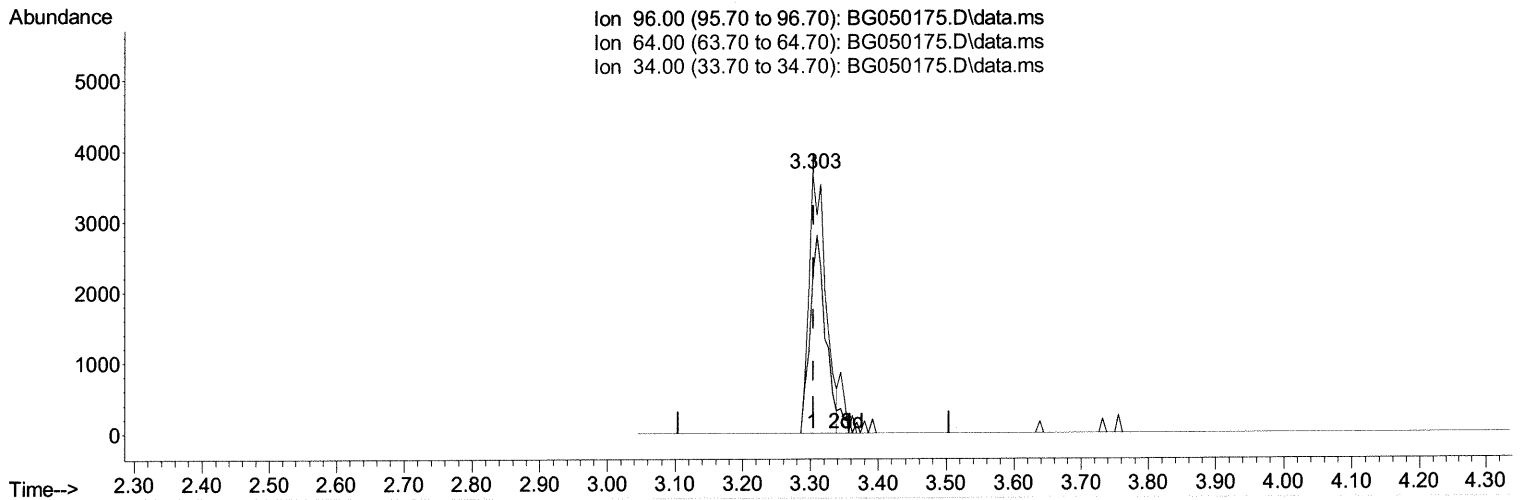
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(3) 1,4-Dioxane-d8 (S)

3.303min (-0.001) 9.28 ng/uL

response 6302

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	58.95
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

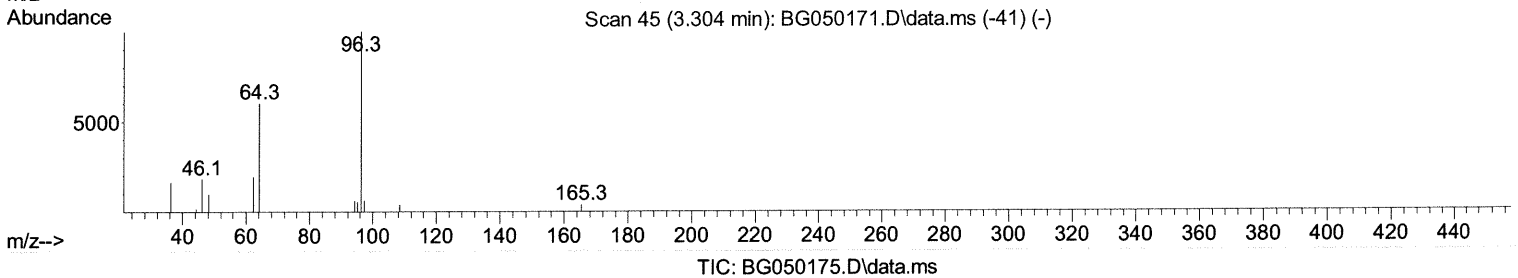
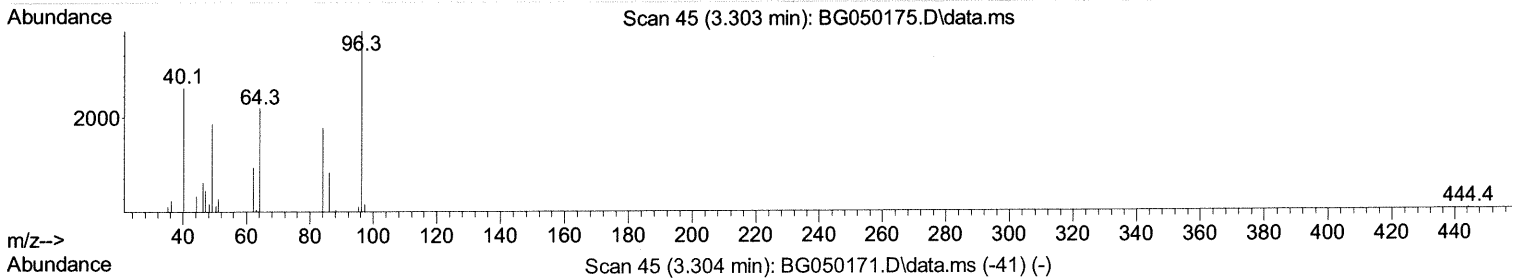
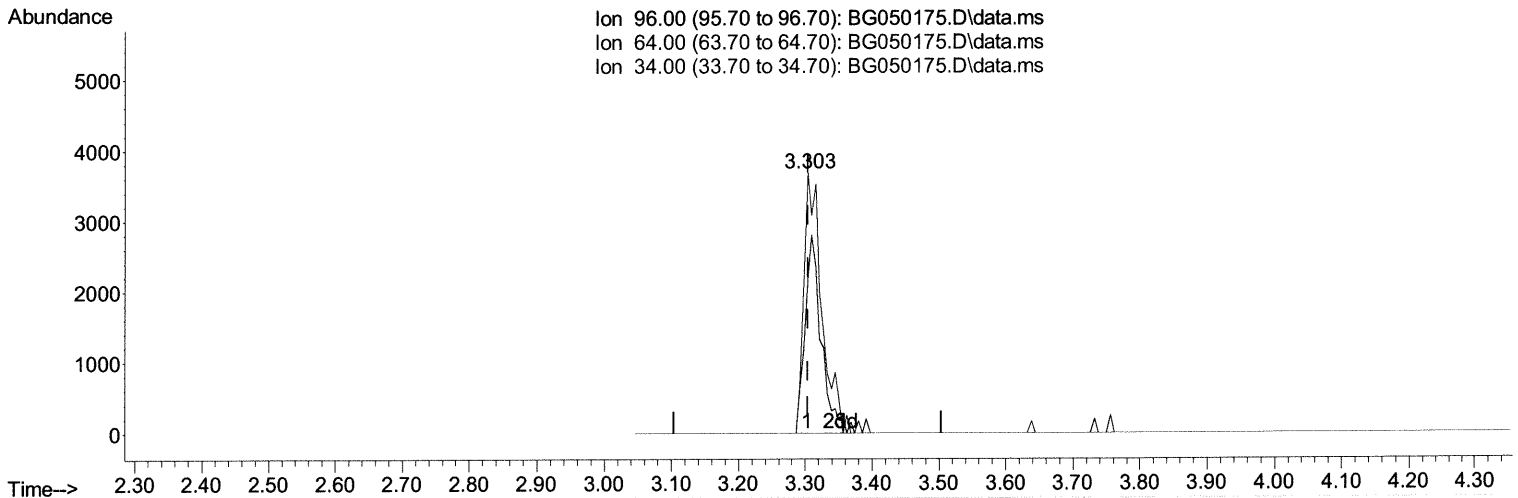
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TIC: BG050175.D\data.ms

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

3.303min (-0.001) 9.98 ng/uL m 09/10/21JU

response 6777

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	58.95
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

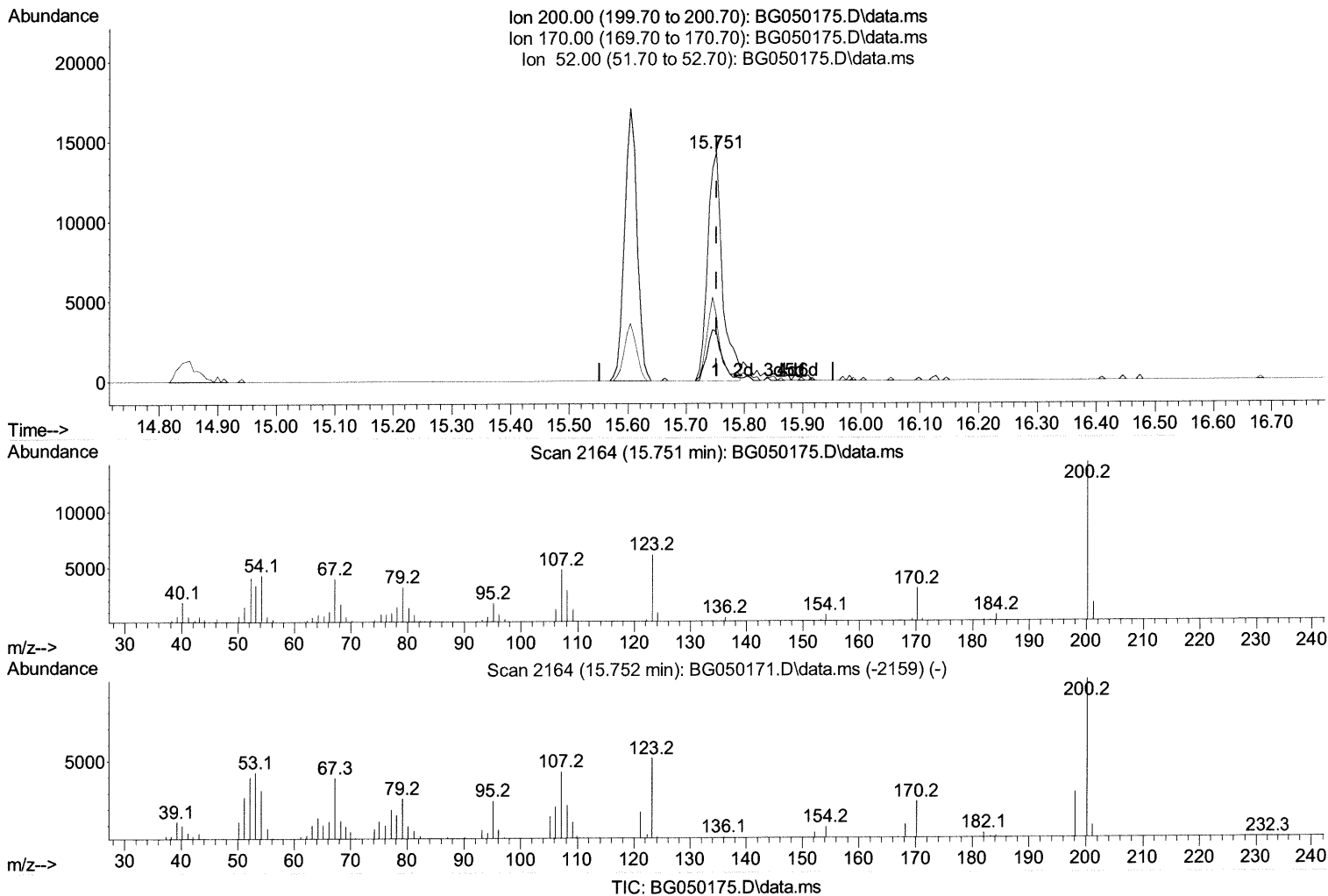
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050175.D
 Acq On : 7 Sep 2021 21:21
 Operator : CG/JU
 Sample : PB138822BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.751min (-0.001) 21.20 ng/ul

response 25397

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.40	21.45
52.00	45.50	28.50#
0.00	0.00	0.00

Quantitation Report (Qedit)

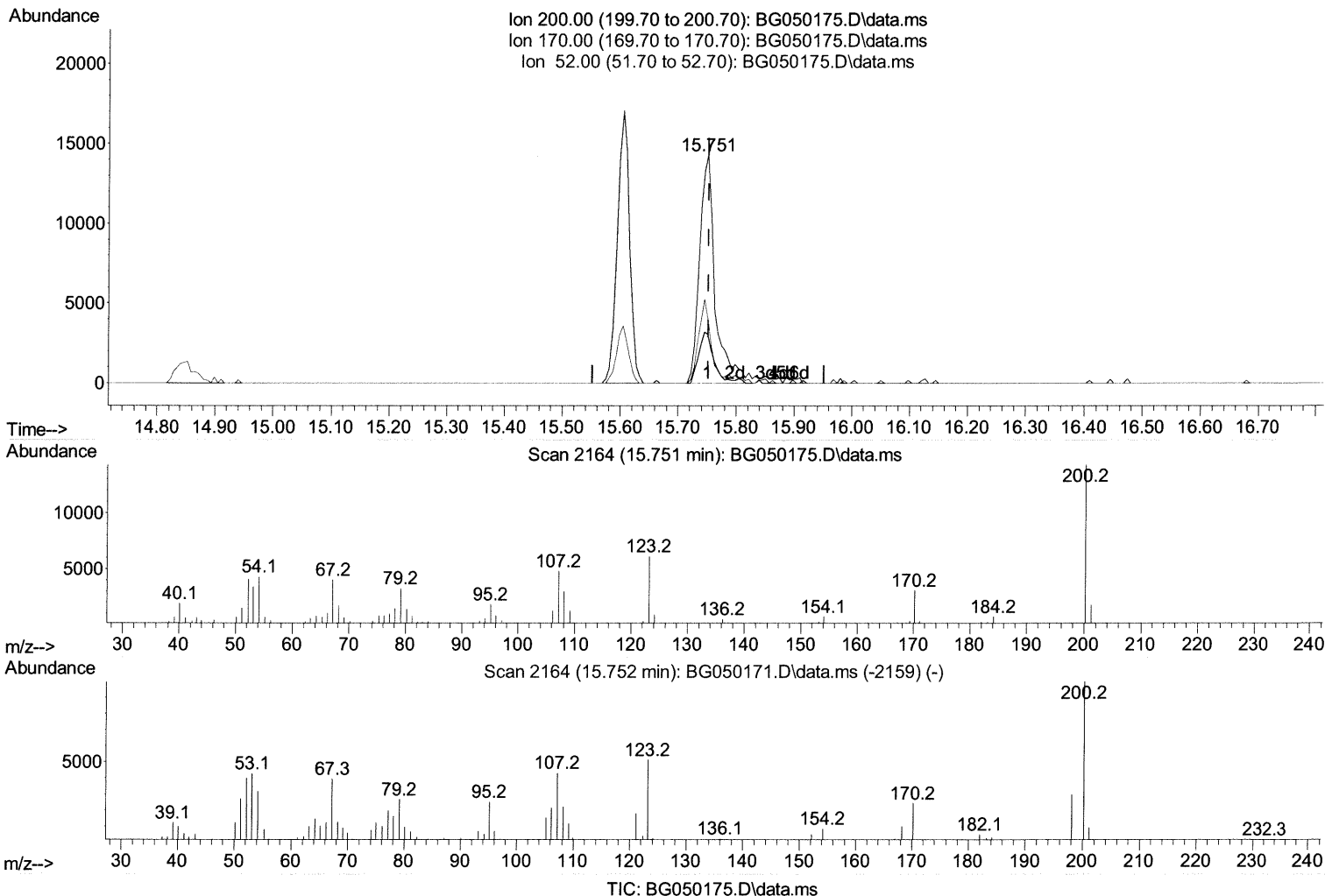
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050175.D
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 Operator : CG/JU
 Sample : PB138822BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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Manual Integrations
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.751min (-0.001) 22.12 ng/ul m 09/10/21 JU

response 26504

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.40	21.45
52.00	45.50	28.50#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	33585	20.000	ng/ul	0.00
20) Naphthalene-d8	10.764	136	146120	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.611	164	89728	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.361	188	198603	20.000	ng/ul	0.00
79) Chrysene-d12	21.631	240	207961	20.000	ng/ul	0.00
88) Perylene-d12	24.845	264	205664	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.303	96	6777m >	9.976	ng/uL >	0.00 09/10/21 JU
4) Pyridine-d5	3.738	84	89292	37.029	ng/ul	0.00
7) Phenol-d5	7.122	99	98337	37.711	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.268	67	55894	36.702	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.480	132	64259	29.692	ng/ul	0.00
15) 4-Methylphenol-d8	8.672	113	81206	39.607	ng/ul	0.00
21) Nitrobenzene-d5	9.119	128	39866	39.749	ng/ul	0.00
24) 2-Nitrophenol-d4	9.847	143	38381	31.471	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	75742	32.540	ng/ul	0.00
31) 4-Chloroaniline-d4	10.911	131	86793	30.266	ng/ul	0.00
46) Dimethylphthalate-d6	14.012	166	217620	32.120	ng/ul	0.00
49) Acenaphthylene-d8	14.300	160	260094	32.240	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.846	143	22799	27.889	ng/ul	0.00
60) Fluorene-d10	15.604	176	198894	35.355	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	26504m >	22.123	ng/ul >	0.00 09/10/21 JU
73) Anthracene-d10	17.460	188	291193	32.661	ng/ul	0.00
81) Pyrene-d10	19.757	212	349402	33.731	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.627	264	344363	31.630	ng/ul	-0.01

Target Compounds Qvalue

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