

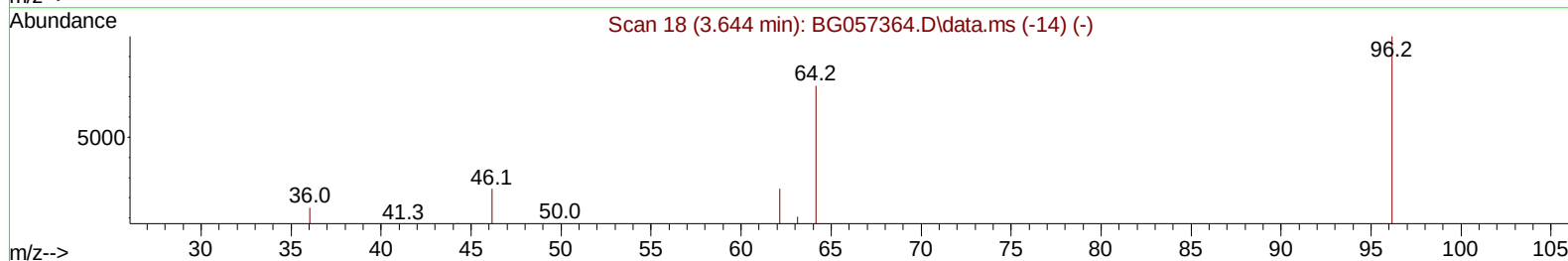
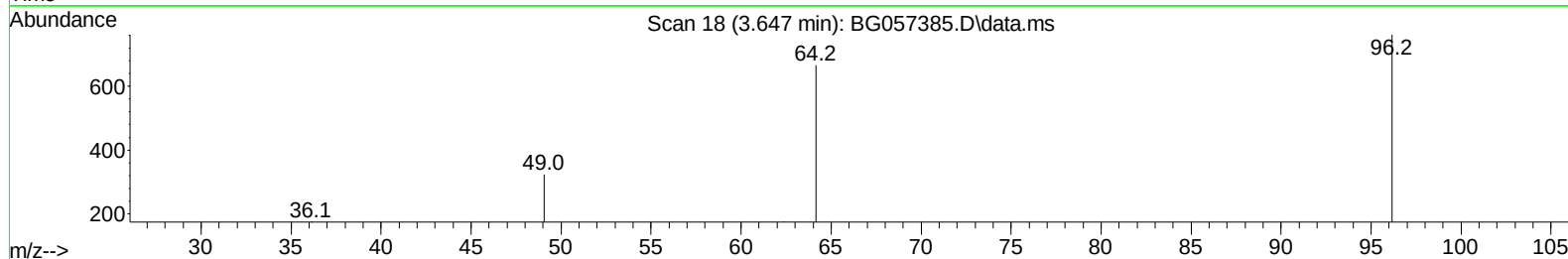
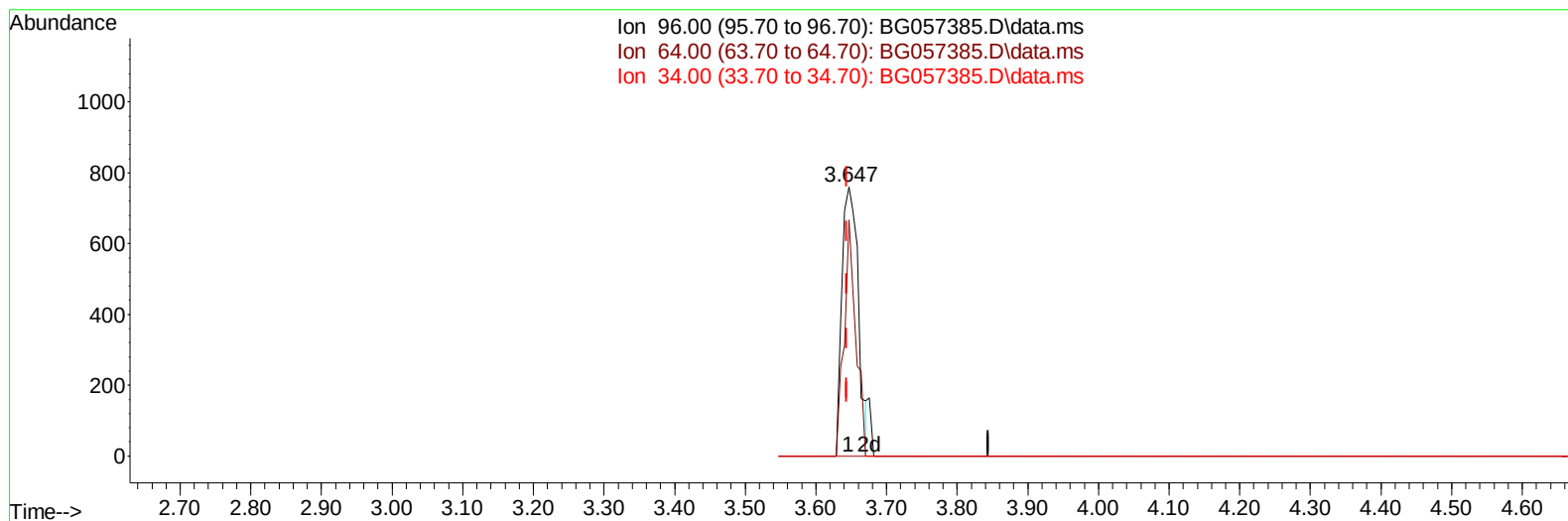
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051023\
Data File : BG057385.D
Acq On : 11 May 2023 1:48
Operator : CG/JU
Sample : 02591-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JREF2

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/11/2023
Supervised By :Jagrut Upadhyay 05/11/2023

Quant Time: May 11 02:45:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 10 11:18:43 2023
Response via : Initial Calibration



TIC: BG057385.D\data.ms

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

3.647min (+ 0.002) 3.27 ng/uL

response 1215

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	70.80	87.52#
34.00	0.00	0.00
0.00	0.00	0.00

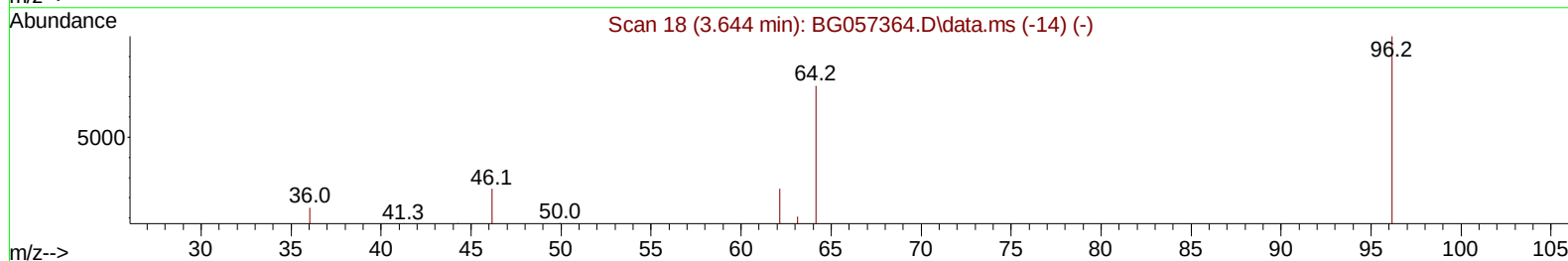
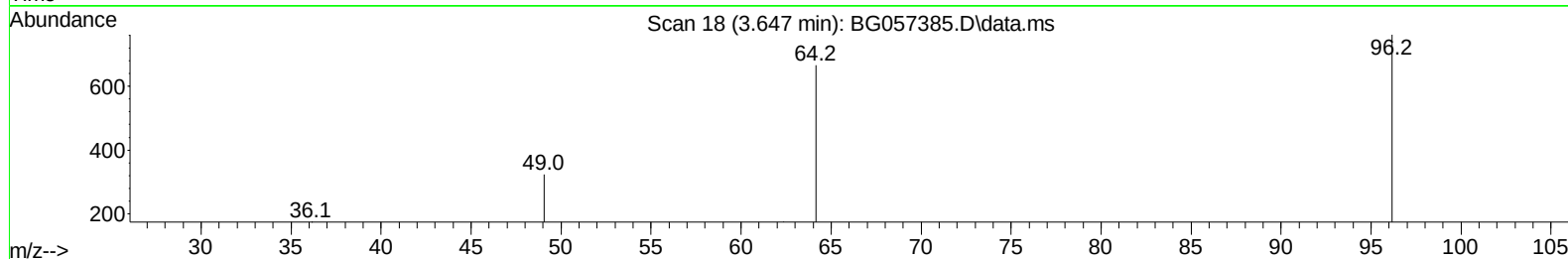
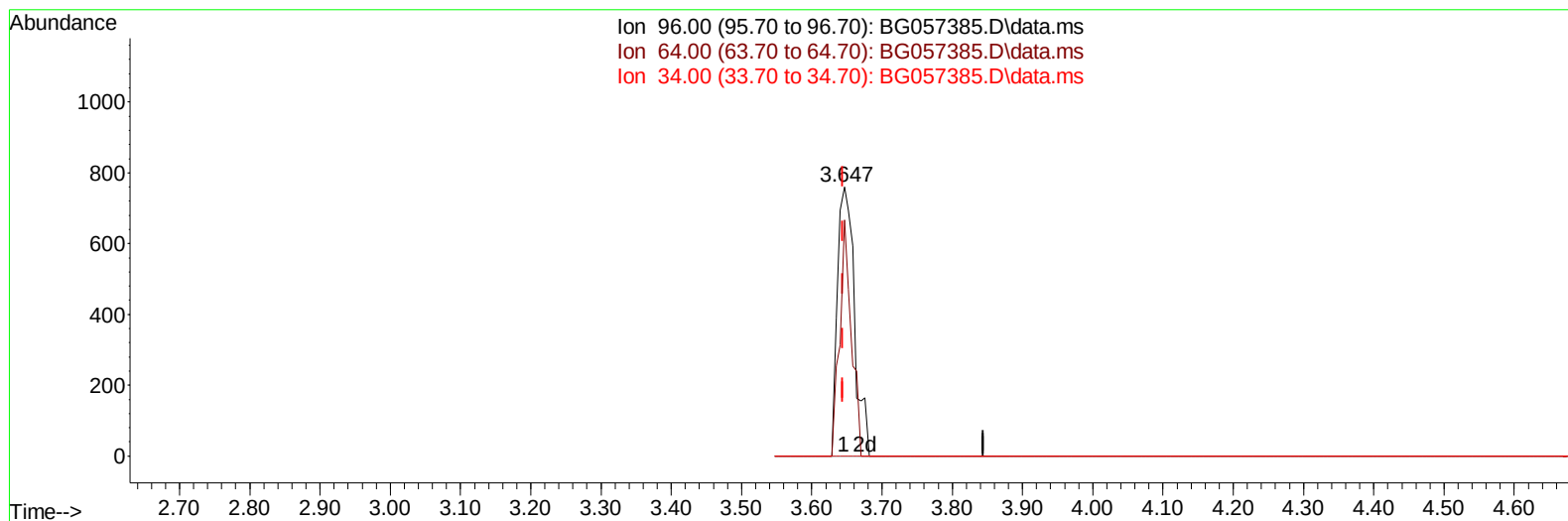
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TIC: BG057385.D\data.ms

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

3.647min (+ 0.002) 3.43 ng/uL m

response 1273

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	70.80	87.52#
34.00	0.00	0.00
0.00	0.00	0.00

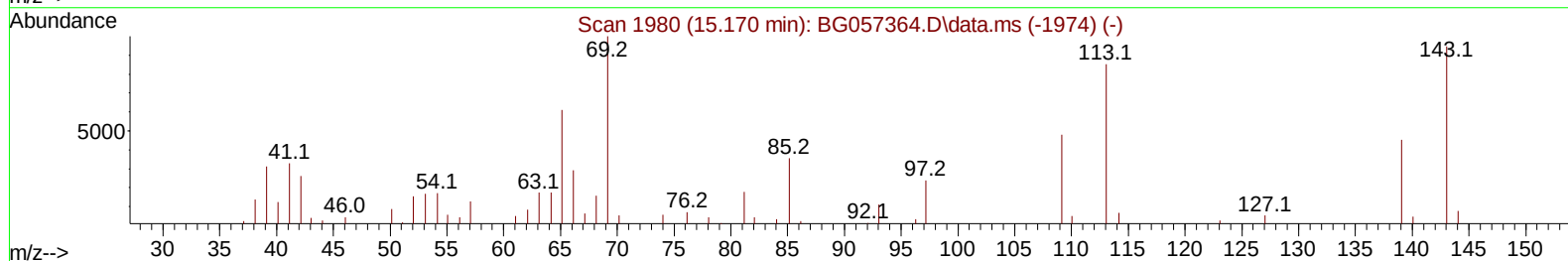
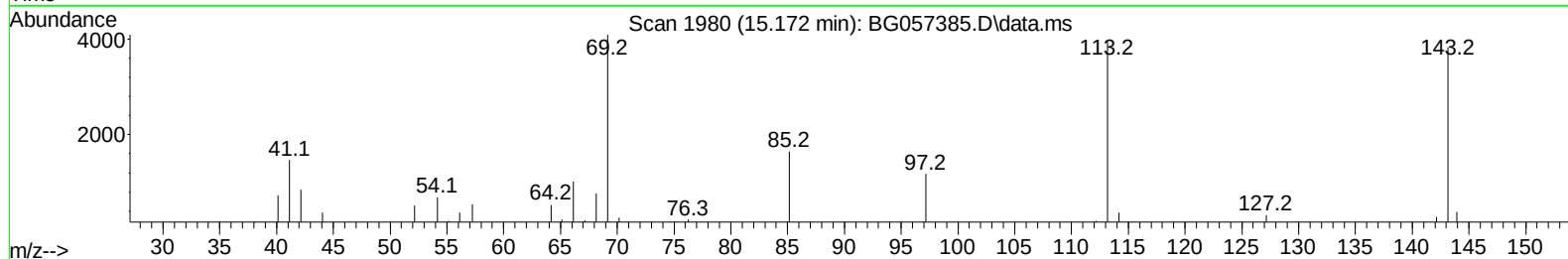
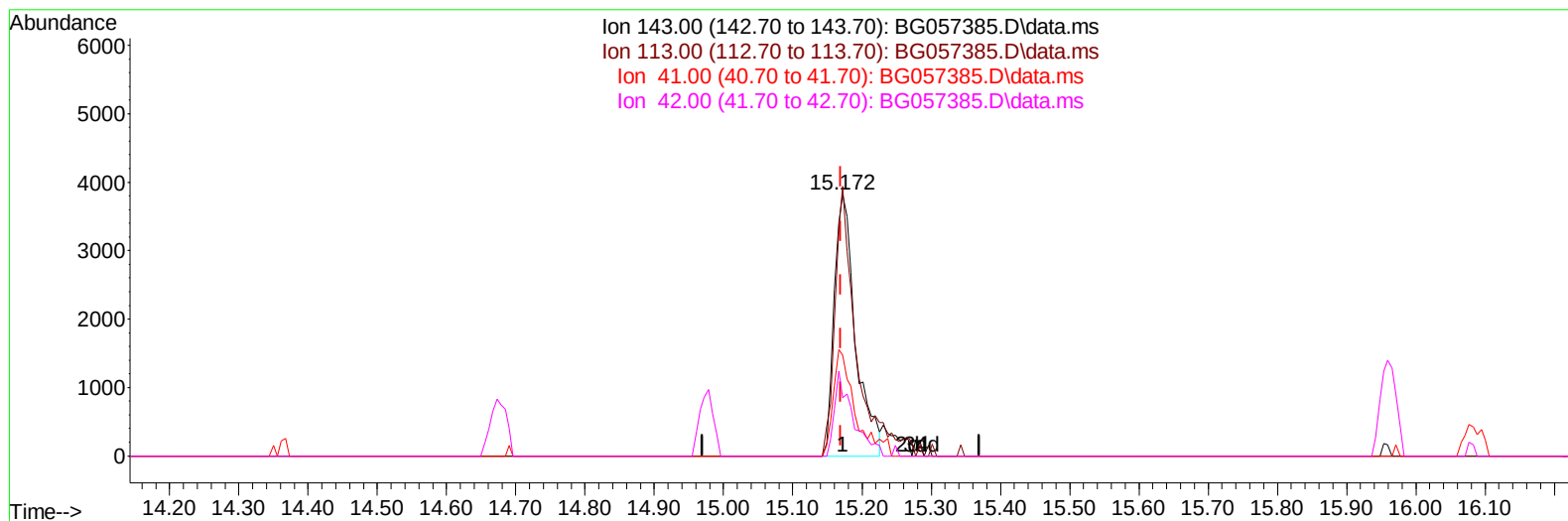
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Manual IntegrationsAPPROVED

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Quant Time: May 11 03:08:25 2023
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Quant Title : SVOA CALIBRATION
QLast Update : Wed May 10 11:18:43 2023
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TIC: BG057385.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.172min (+ 0.002) 12.59 ng/ul

response 8158

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	103.01
41.00	44.40	38.49
42.00	28.40	22.24#

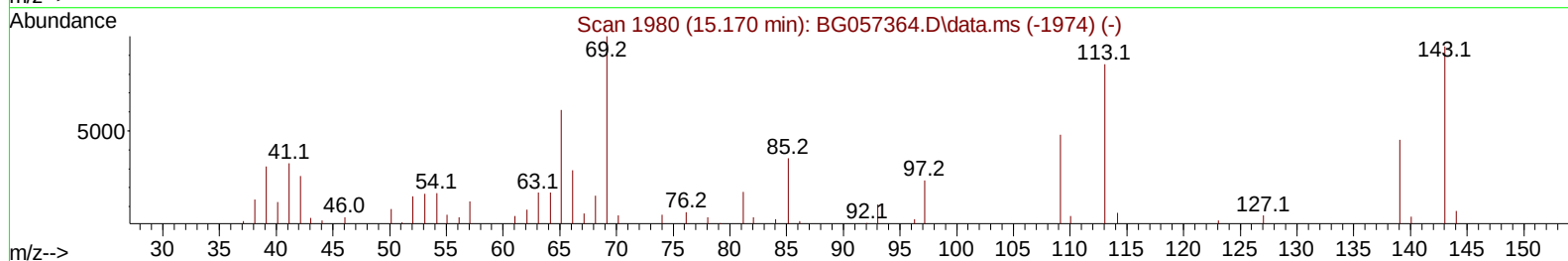
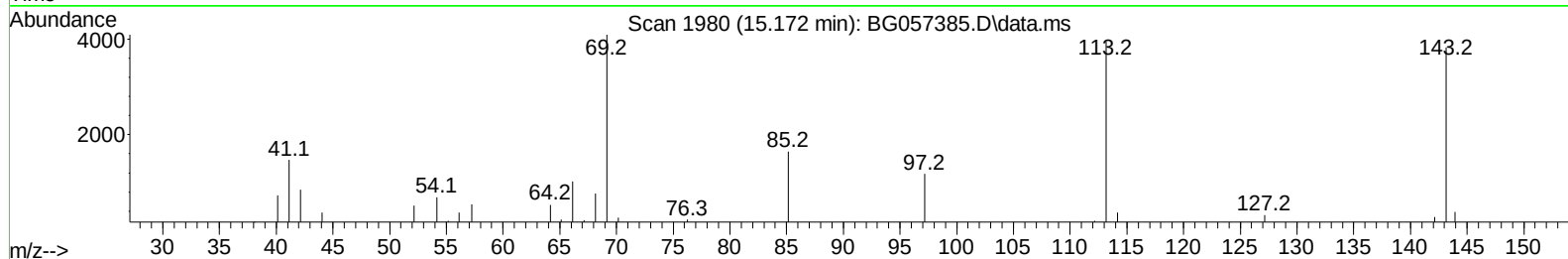
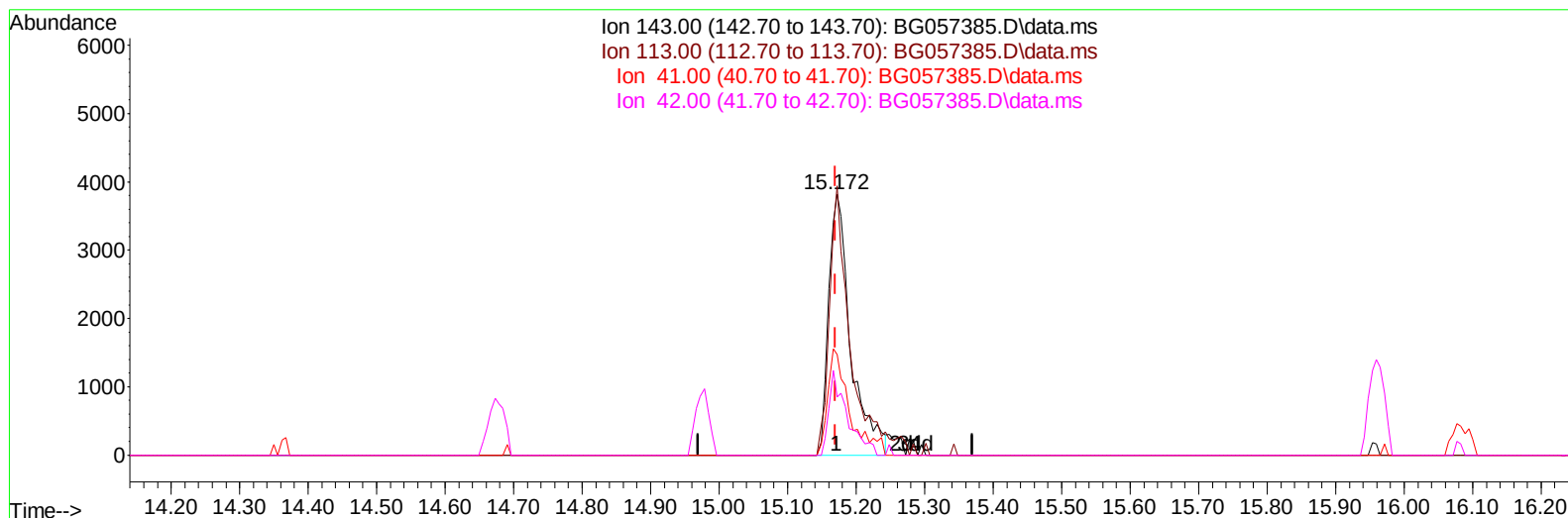
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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG057385.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.172min (+ 0.002) 13.18 ng/ul m

response 8537

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	103.01
41.00	44.40	38.49
42.00	28.40	22.24#

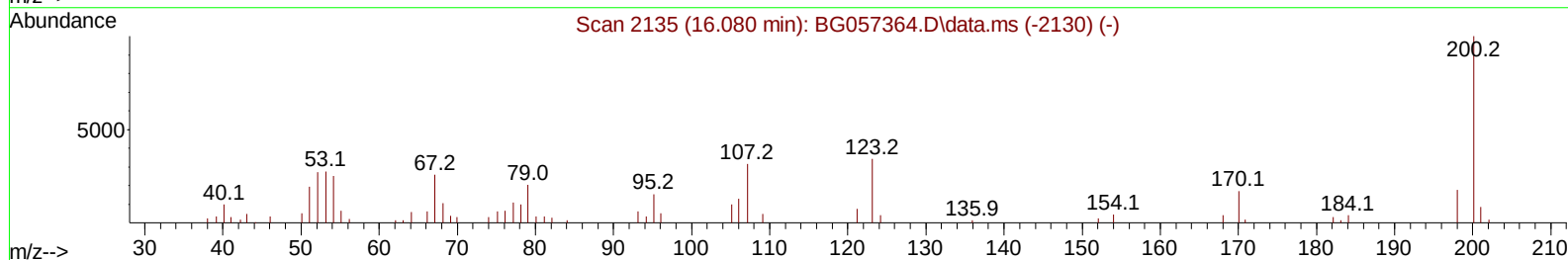
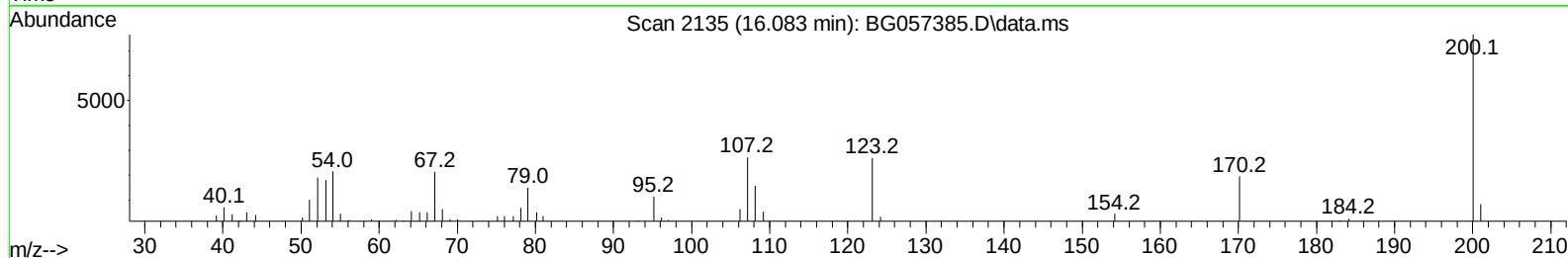
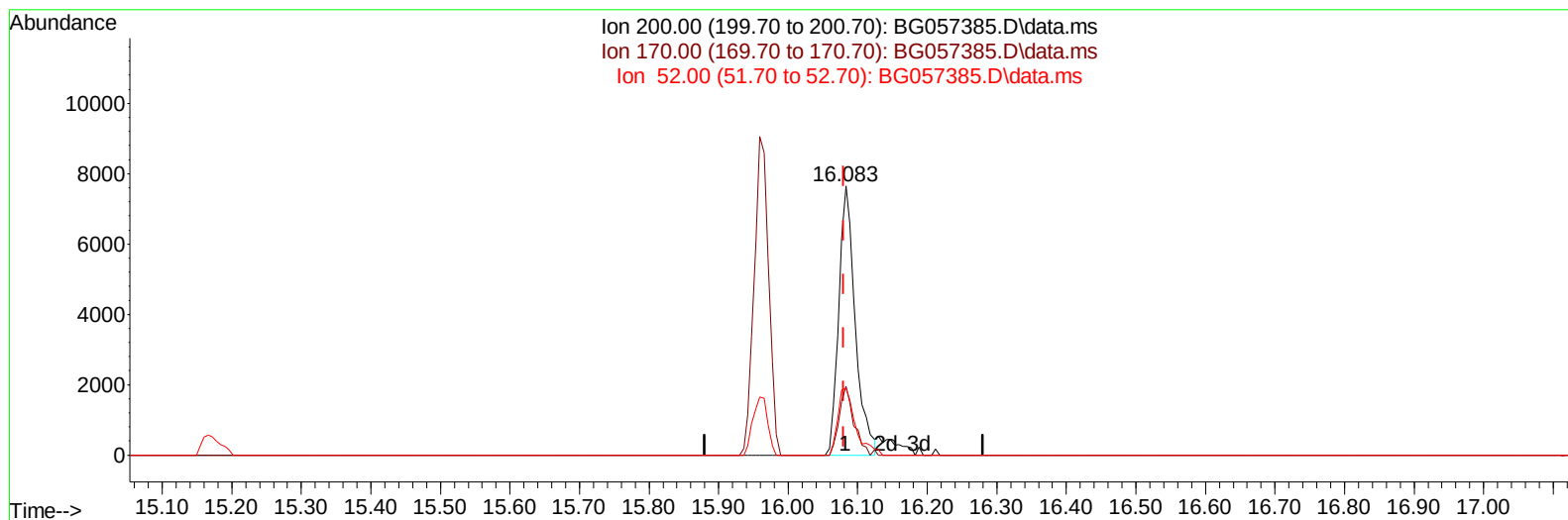
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 Acq On : 11 May 2023 1:48
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 Sample : 02591-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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TIC: BG057385.D\data.ms

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

16.083min (+ 0.003) 15.74 ng/ul

response 12626

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.80	25.60
52.00	34.60	24.86#
0.00	0.00	0.00

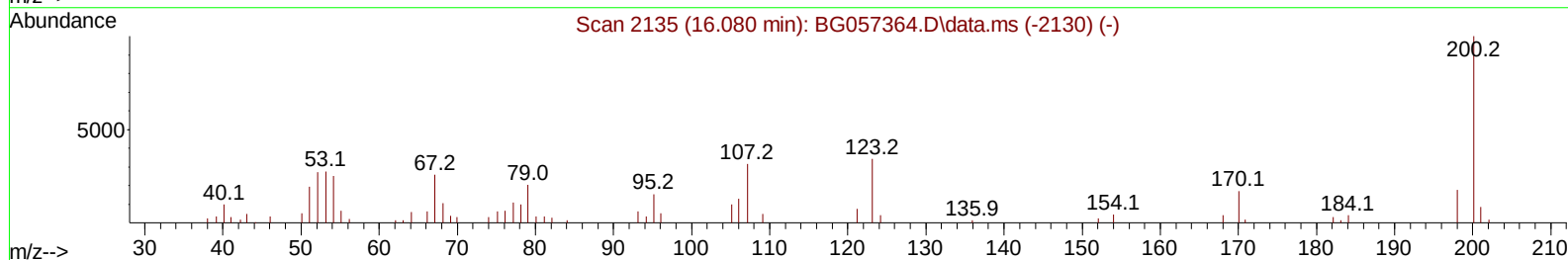
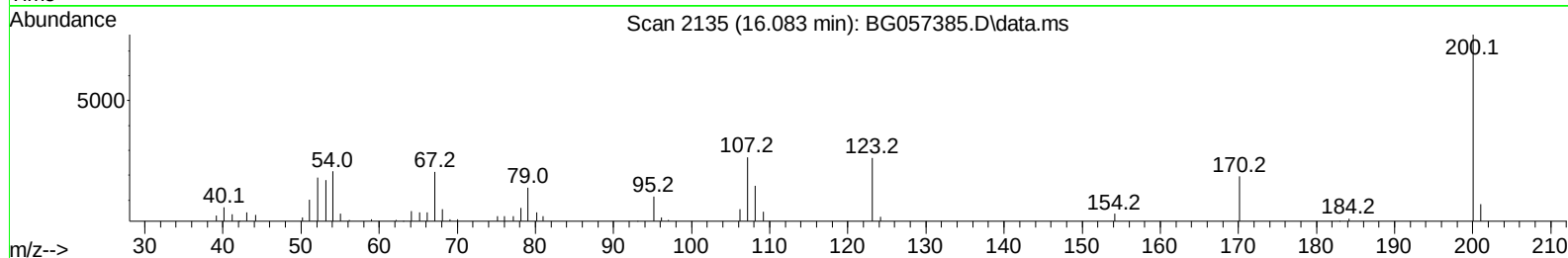
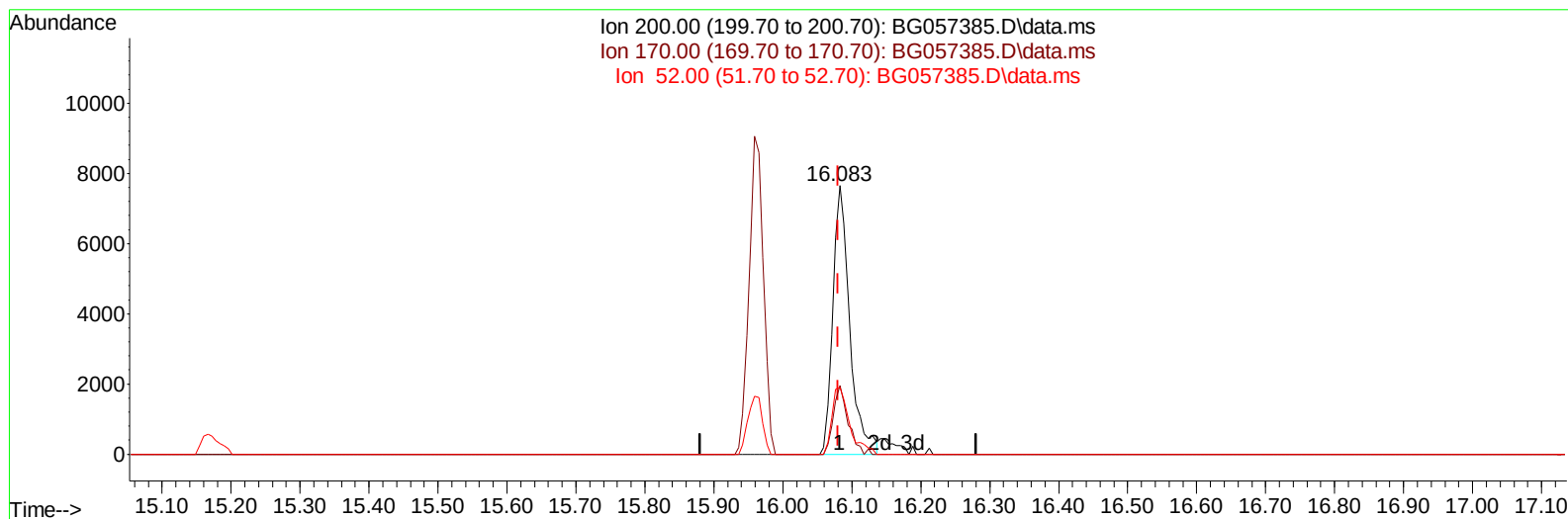
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 Operator : CG/JU
 Sample : 02591-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BG057385.D\data.ms

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

16.083min (+ 0.003) 16.13 ng/ul m

response 12937

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.80	25.60
52.00	34.60	24.86#
0.00	0.00	0.00

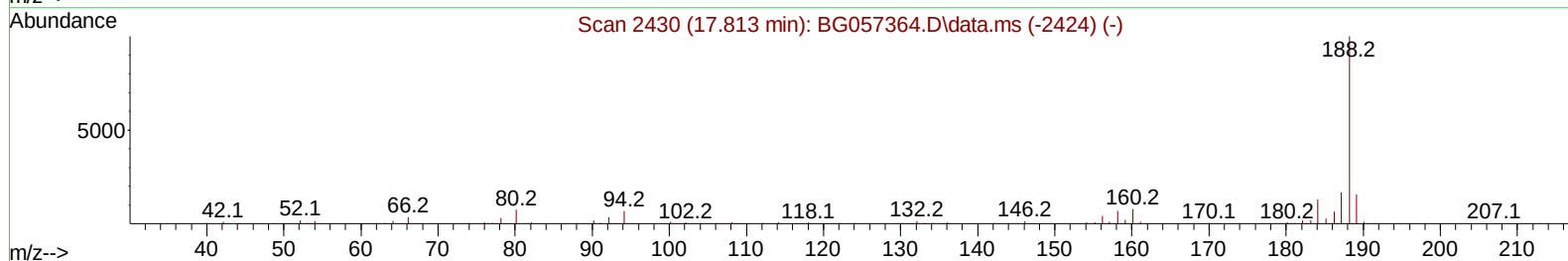
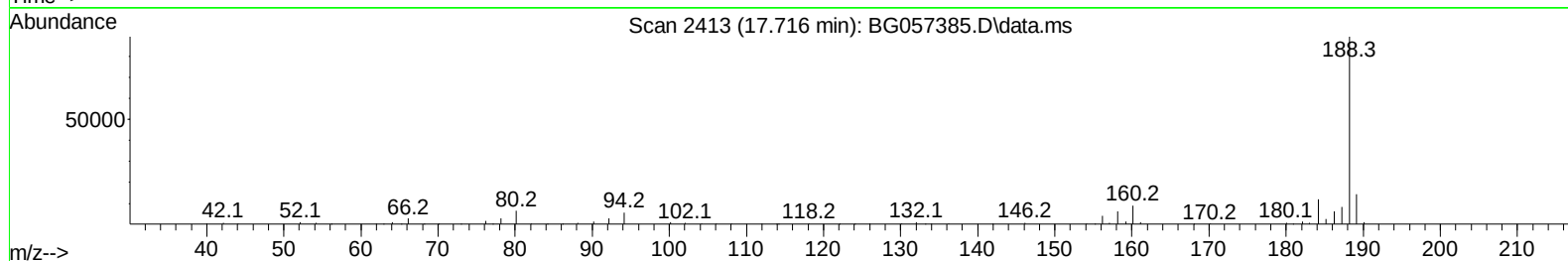
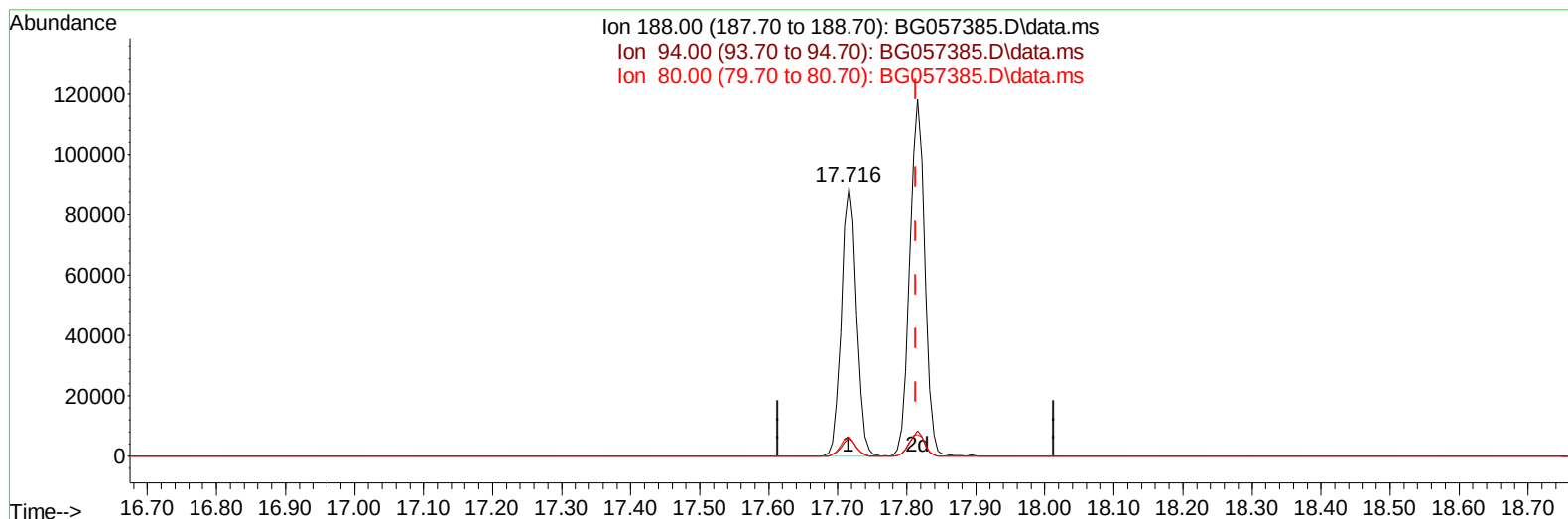
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Instrument :
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TIC: BG057385.D\data.ms

(73) Anthracene-d10 (S)

17.716min (-0.097) 21.91 ng/u1

response 135352

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.90	6.44
80.00	7.70	7.19
0.00	0.00	0.00

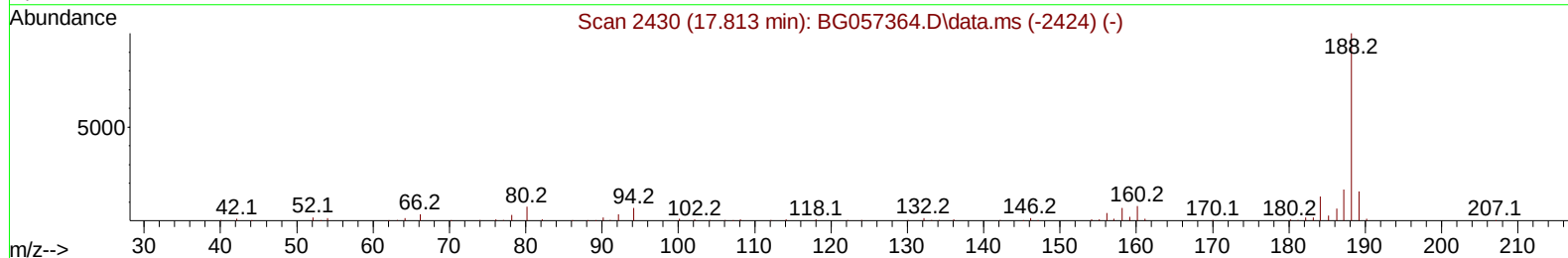
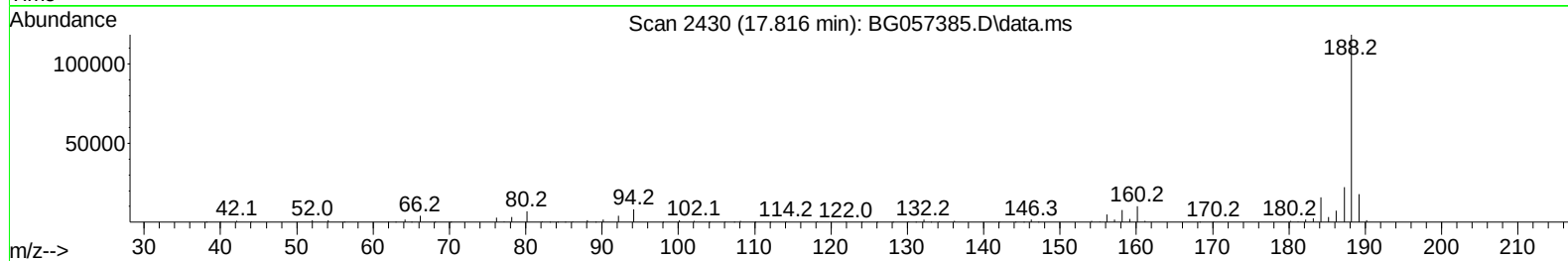
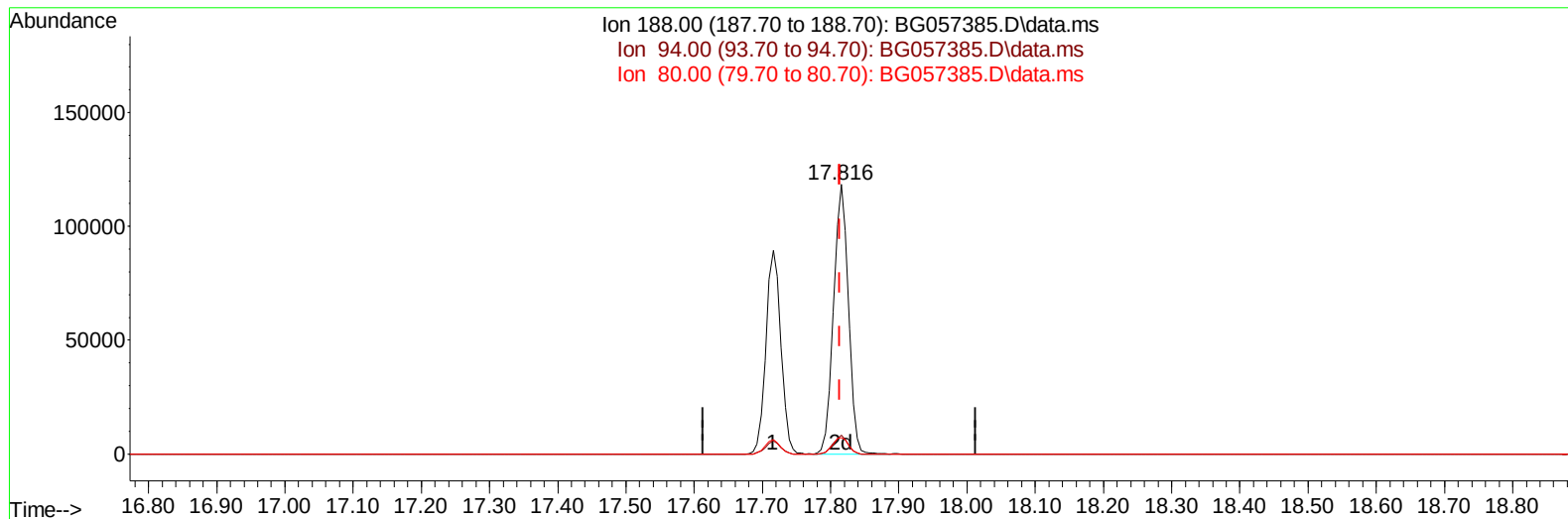
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051023\
 Data File : BG057385.D
 Acq On : 11 May 2023 1:48
 Operator : CG/JU
 Sample : 02591-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/11/2023
 Supervised By :Jagrut Upadhyay 05/11/2023

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 Response via : Initial Calibration



TIC: BG057385.D\data.ms

(73) Anthracene-d10 (S)

17.816min (+ 0.002) 28.90 ng/ul m

response 178520

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.90	7.01
80.00	7.70	6.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051023\
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 Sample : 02591-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/11/2023
 Supervised By :Jagrut Upadhyay 05/11/2023

Quant Time: May 11 03:09:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.364	152	16344	20.000	ng/ul	0.00
20) Naphthalene-d8	11.201	136	69597	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.978	164	56176	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.716	188	135352	20.000	ng/ul	0.00
79) Chrysene-d12	22.028	240	124576	20.000	ng/ul	0.00
88) Perylene-d12	25.529	264	131149	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.647	96	1273m	3.427	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.506	99	32375	20.946	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.671	67	19356	21.291	ng/ul	0.00
11) 2-Chlorophenol-d4	7.894	132	24382	23.937	ng/ul	0.00
15) 4-Methylphenol-d8	9.063	113	24430	20.796	ng/ul	0.00
21) Nitrobenzene-d5	9.544	128	13348	24.083	ng/ul	0.00
24) 2-Nitrophenol-d4	10.273	143	16681	25.789	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.819	165	32675	26.183	ng/ul	0.00
31) 4-Chloroaniline-d4	11.330	131	6927	4.188	ng/ul	0.00
46) Dimethylphthalate-d6	14.362	166	114153	25.645	ng/ul	0.00
49) Acenaphthylene-d8	14.679	160	136815	27.347	ng/ul	0.00
54) 4-Nitrophenol-d4	15.172	143	8537m	13.176	ng/ul	0.00
60) Fluorene-d10	15.959	176	112580	27.150	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.083	200	12937m	16.132	ng/ul	0.00
73) Anthracene-d10	17.816	188	178520m	28.895	ng/ul	0.00
81) Pyrene-d10	20.083	212	213788	28.912	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.288	264	176133	26.361	ng/ul	0.00

Target Compounds Qvalue

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

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