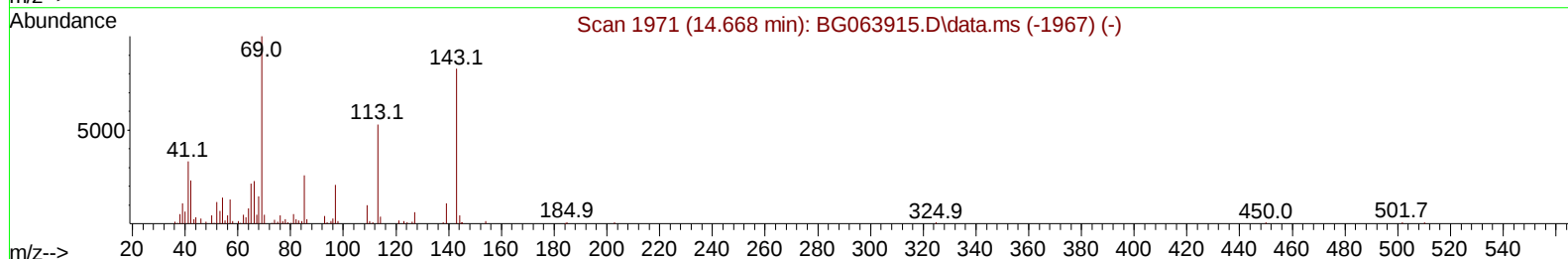
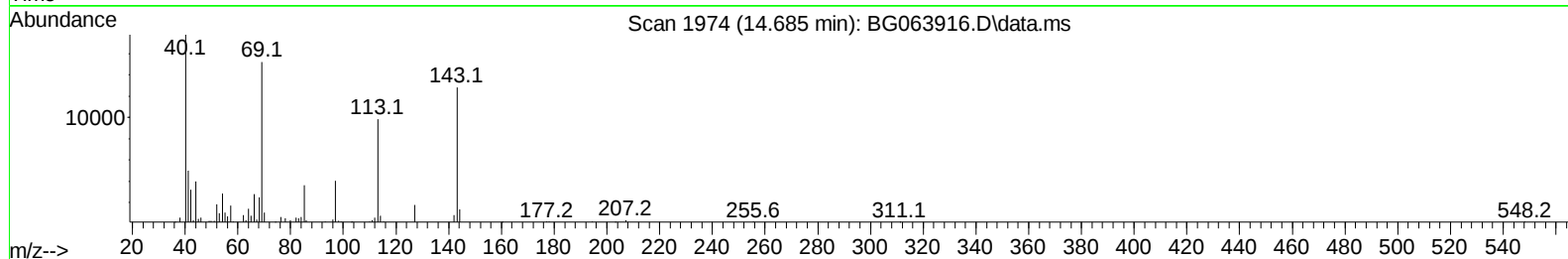
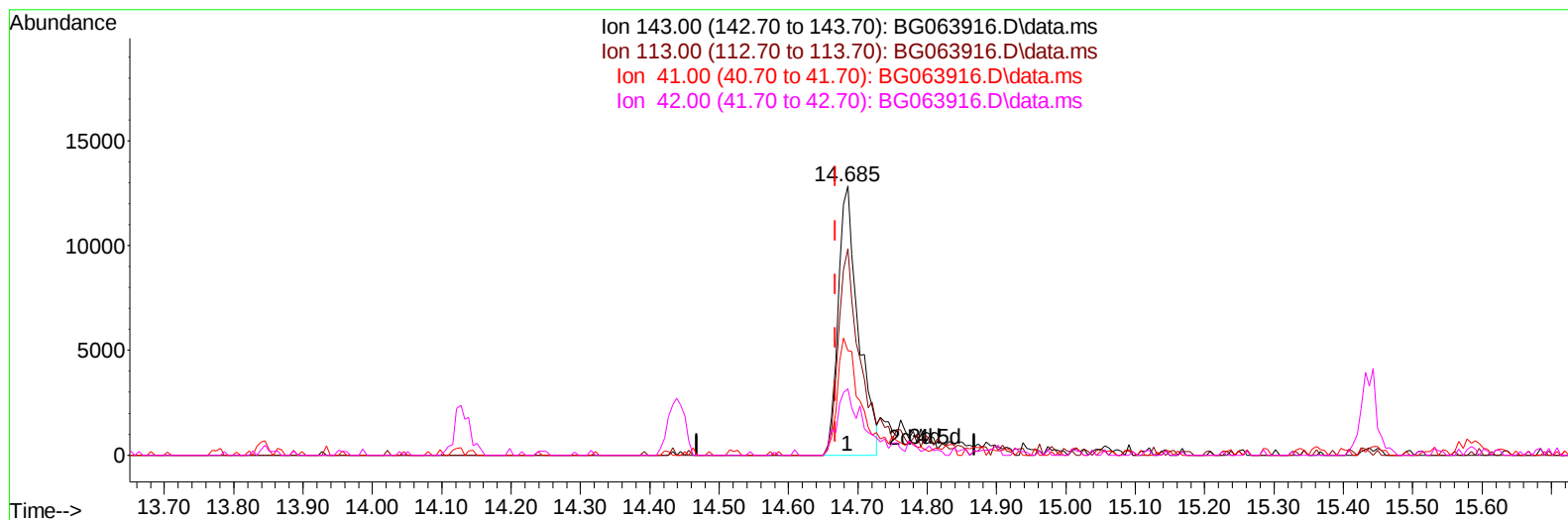


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123124\
Data File : BG063916.D
Acq On : 31 Dec 2024 10:50
Operator : RC/JU
Sample : PB165853BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 01 00:45:47 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 24 00:44:39 2024
Response via : Initial Calibration



TIC: BG063916.D\data.ms

(54) 4-Nitrophenol-d4 (S)

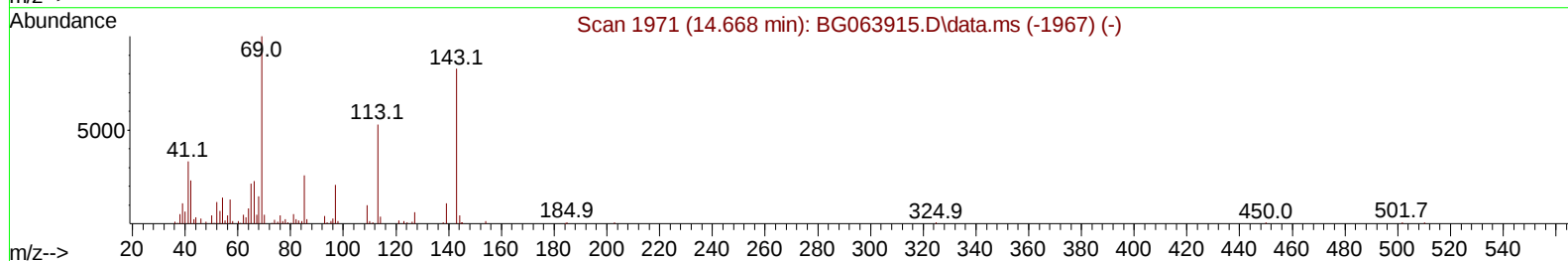
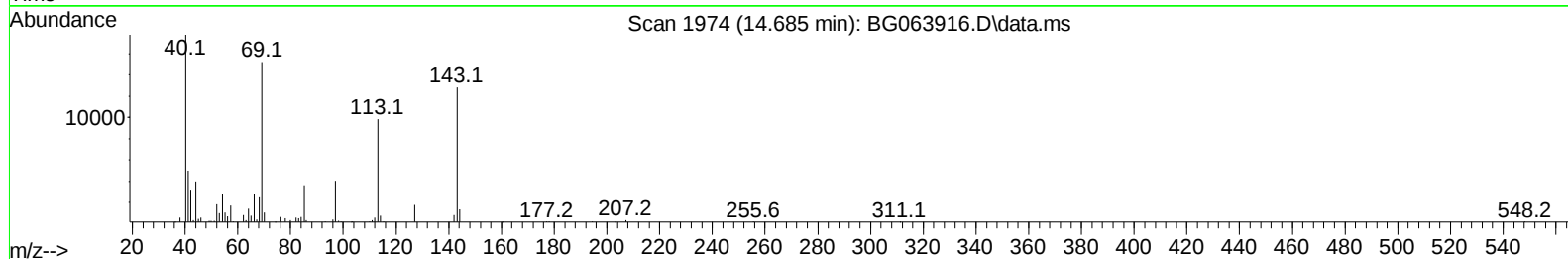
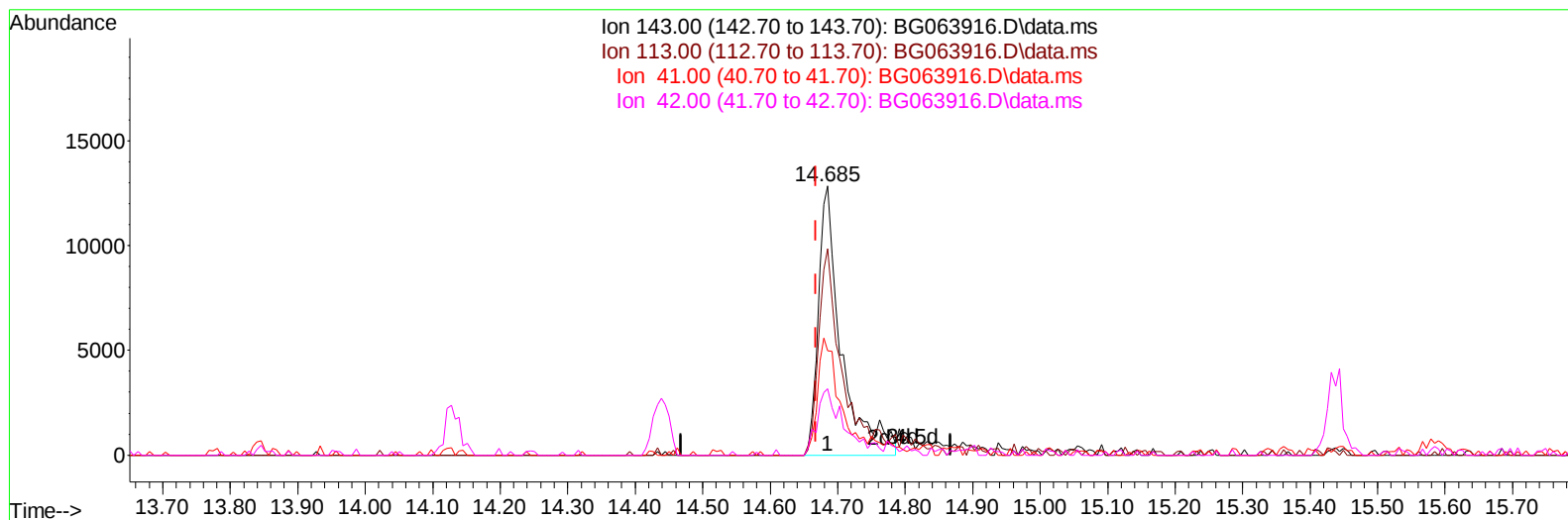
14.685min (+ 0.017) 16.29 ng/u1

response 25833

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	89.70	76.69
41.00	48.20	38.90
42.00	26.60	24.75

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123124\
Data File : BG063916.D
Acq On : 31 Dec 2024 10:50
Operator : RC/JU
Sample : PB165853BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 01 00:45:47 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 24 00:44:39 2024
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TIC: BG063916.D\data.ms

(54) 4-Nitrophenol-d4 (S)

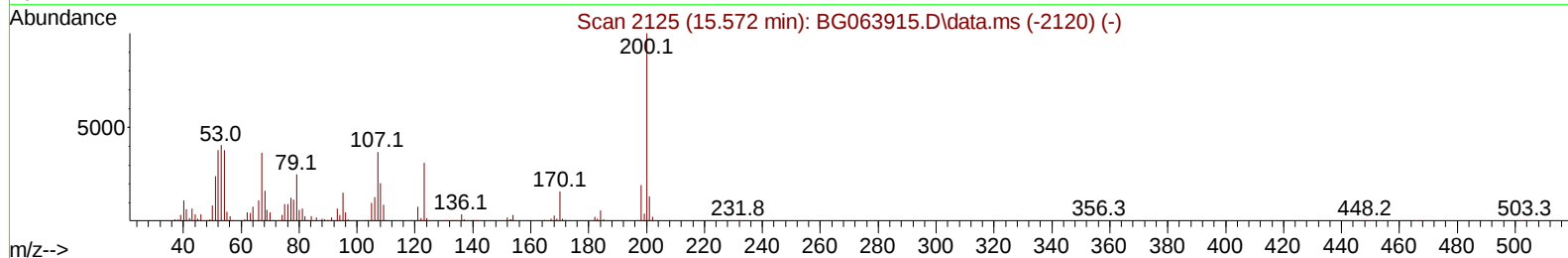
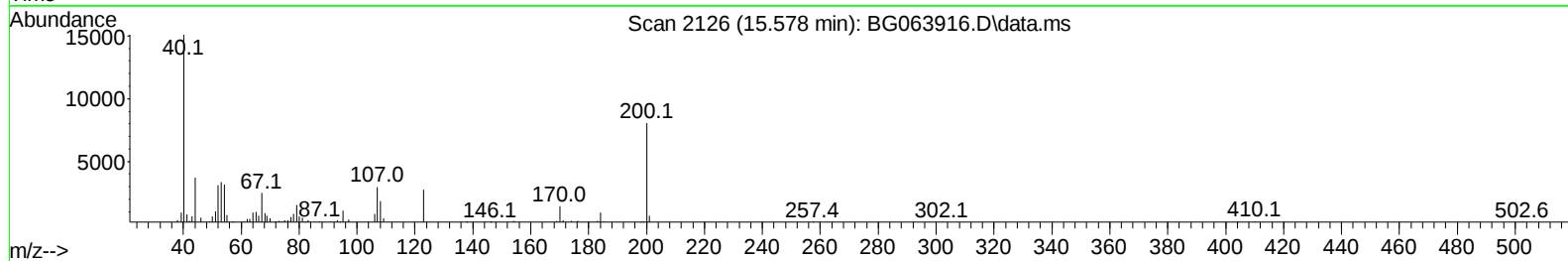
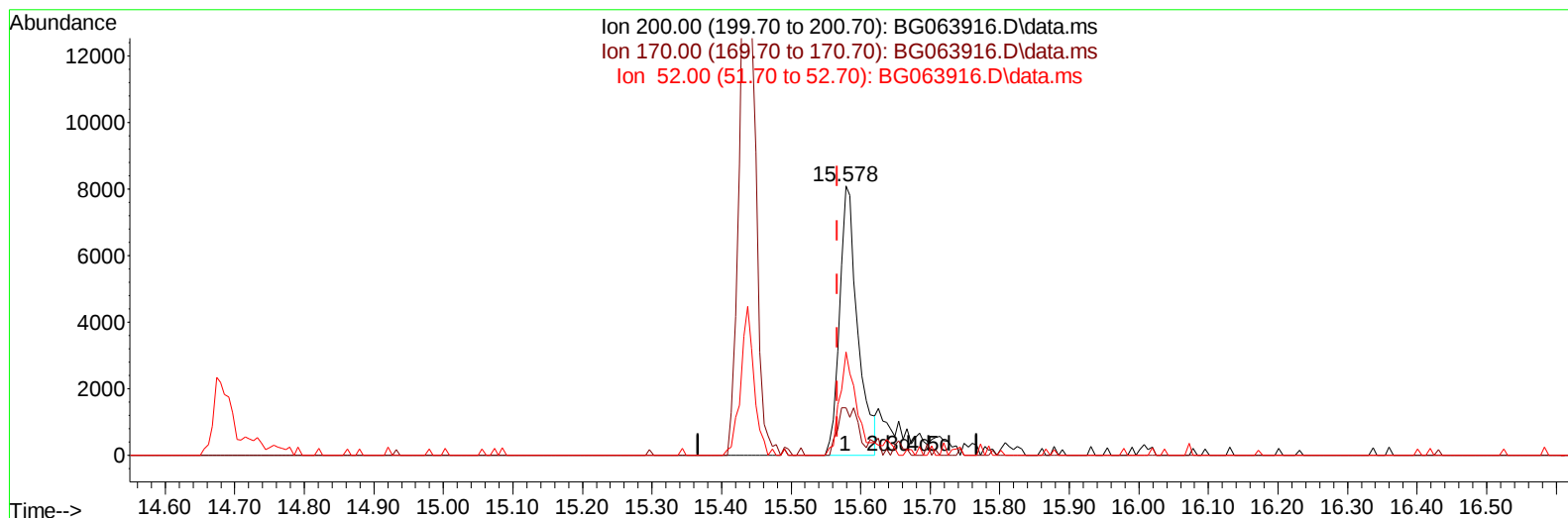
14.685min (+ 0.017) 19.03 ng/ul m

response 30176

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	89.70	76.69
41.00	48.20	38.90
42.00	26.60	24.75

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123124\
Data File : BG063916.D
Acq On : 31 Dec 2024 10:50
Operator : RC/JU
Sample : PB165853BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 01 00:45:47 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 24 00:44:39 2024
Response via : Initial Calibration



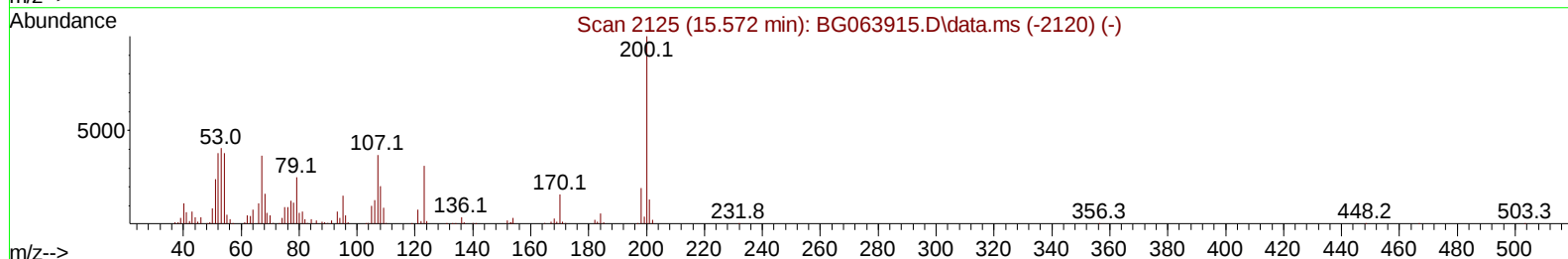
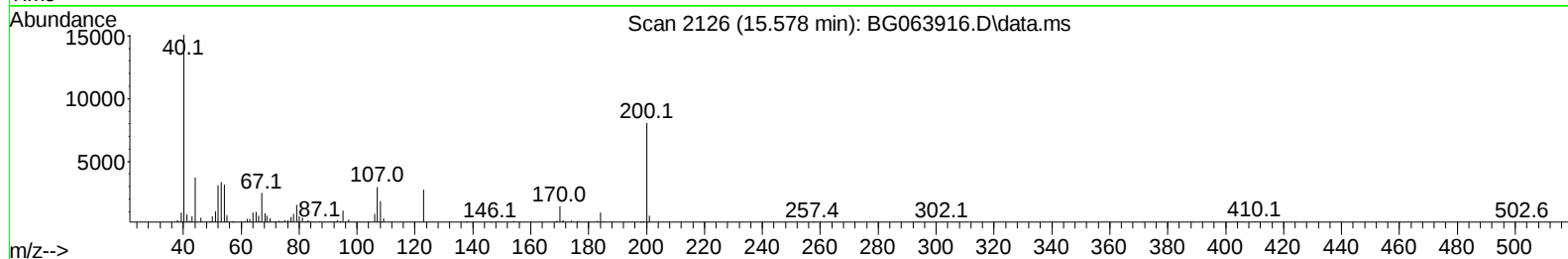
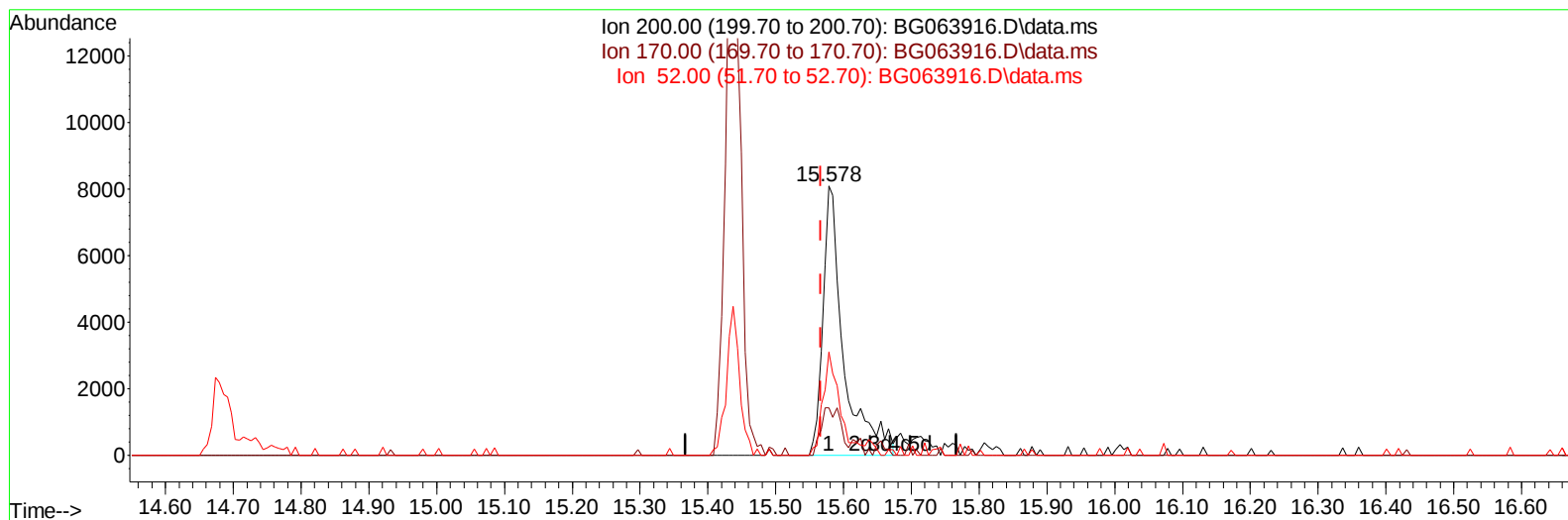
TIC: BG063916.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.578min (+ 0.011) 7.11 ng/ul****response 14616**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.90	17.48#
52.00	43.10	38.43
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123124\
Data File : BG063916.D
Acq On : 31 Dec 2024 10:50
Operator : RC/JU
Sample : PB165853BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 01 00:45:47 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 24 00:44:39 2024
Response via : Initial Calibration



TIC: BG063916.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.578min (+ 0.011) 8.38 ng/ul m****response 17219**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.90	17.48#
52.00	43.10	38.43
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123124\
 Data File : BG063916.D
 Acq On : 31 Dec 2024 10:50
 Operator : RC/JU
 Sample : PB165853BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 01 01:19:29 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	62997	20.000	ng/ul	0.00
20) Naphthalene-d8	10.584	136	246053	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	164441	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.194	188	395984	20.000	ng/ul	0.00
79) Chrysene-d12	21.454	240	463273	20.000	ng/ul	0.02
88) Perylene-d12	24.474	264	486569	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	8833	5.495	ng/uL	0.00
4) Pyridine-d5	3.669	84	128991	25.436	ng/ul	0.00
7) Phenol-d5	6.983	99	130755	22.536	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.124	67	91349	22.796	ng/ul	0.00
11) 2-Chlorophenol-d4	7.335	132	86114	22.926	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	96108	21.336	ng/ul	0.00
21) Nitrobenzene-d5	8.951	128	42372	24.118	ng/ul	0.00
24) 2-Nitrophenol-d4	9.668	143	46749	23.739	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.220	165	87003	22.819	ng/ul	0.00
31) 4-Chloroaniline-d4	10.725	131	105573	21.807	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	273173	24.413	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	326305	25.582	ng/ul	0.00
54) 4-Nitrophenol-d4	14.685	143	30176m	19.034	ng/ul	0.02
60) Fluorene-d10	15.437	176	236881	23.944	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.578	200	17219m	8.375	ng/ul	0.01
73) Anthracene-d10	17.294	188	416335	24.455	ng/ul	0.00
81) Pyrene-d10	19.597	212	544114	22.570	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.262	264	556628	24.361	ng/ul	0.02

Target Compounds	Qvalue

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

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