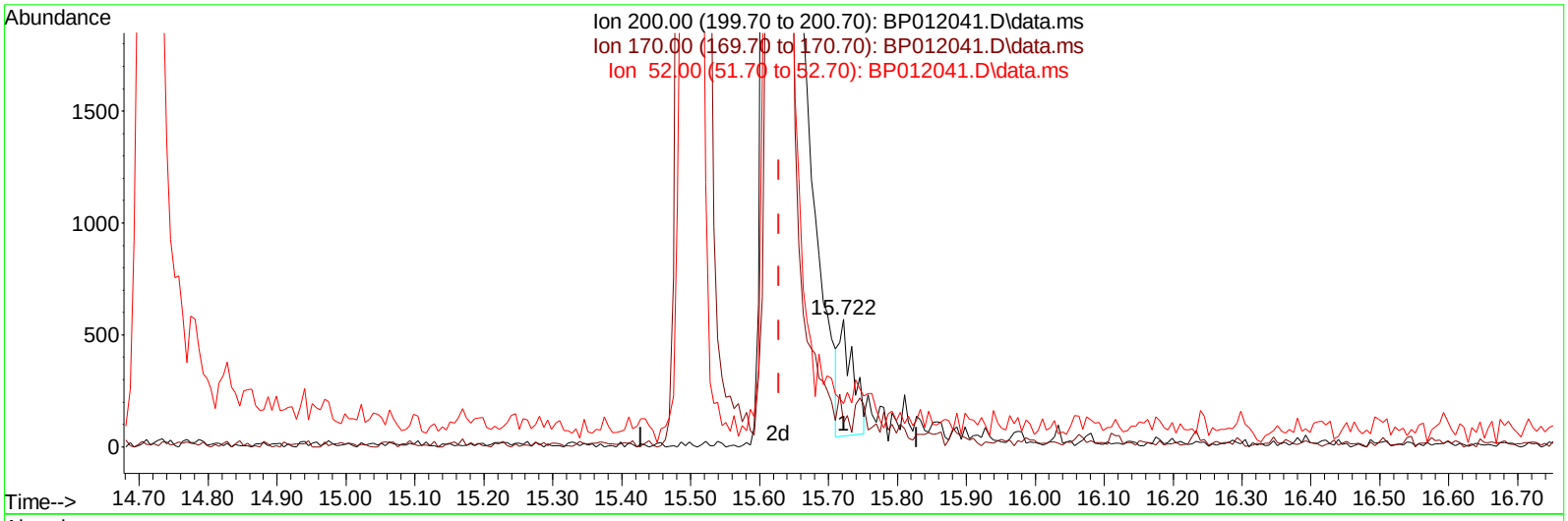
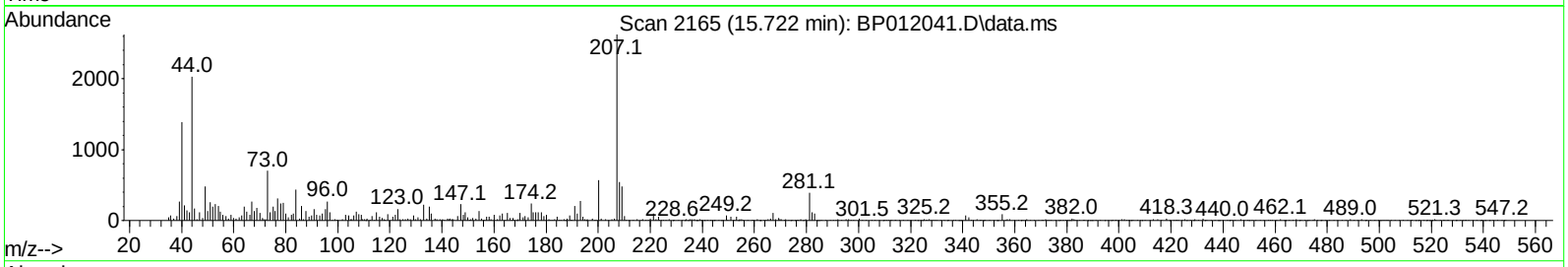
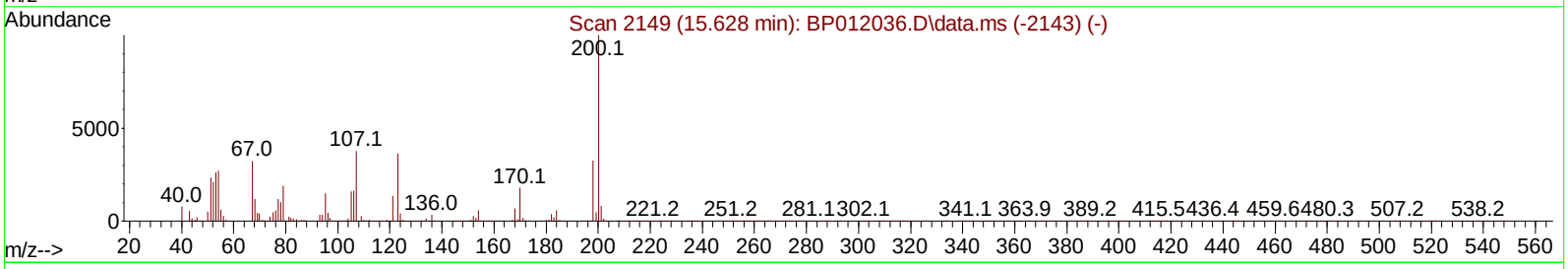
Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012041.D
Acq On : 11 Oct 2022 15:38
Operator : CG/JU
Sample : PB148209BL
Misc :
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Oct 12 00:22:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 11 23:54:43 2022
Response via : Initial Calibration







TIC: BP012041.D\data.ms

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

15.722min (+ 0.094) 0.14 ng/ul

response 748

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	17.90	19.33
52.00	34.50	33.74
0.00	0.00	0.00

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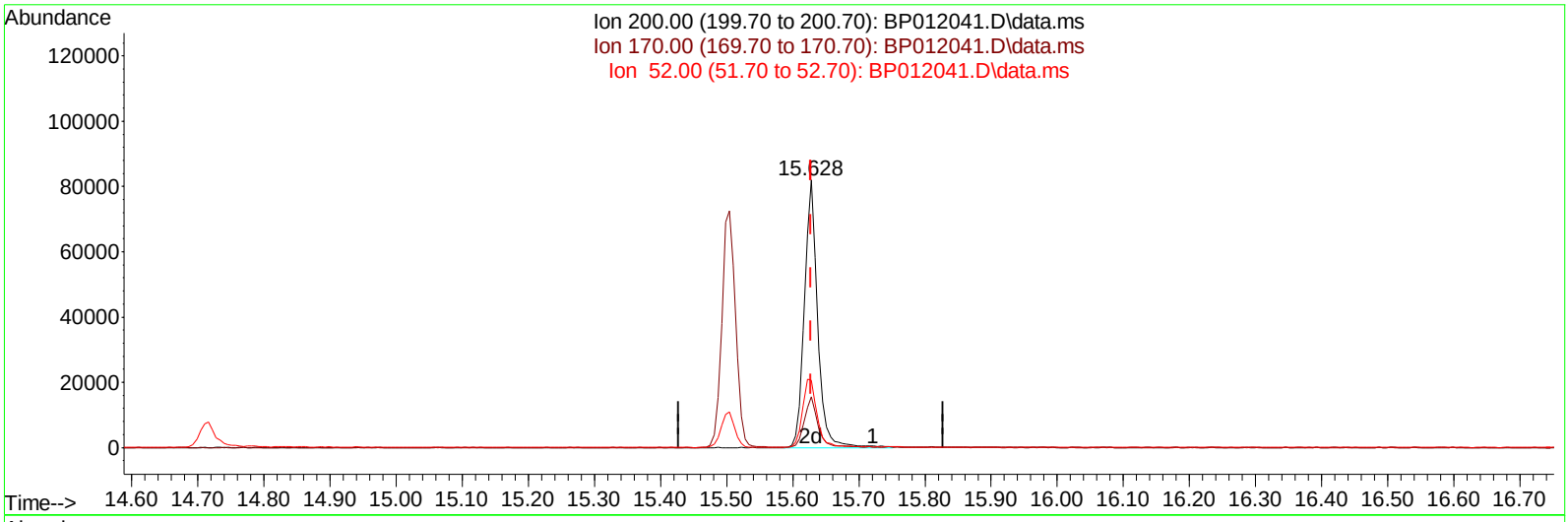

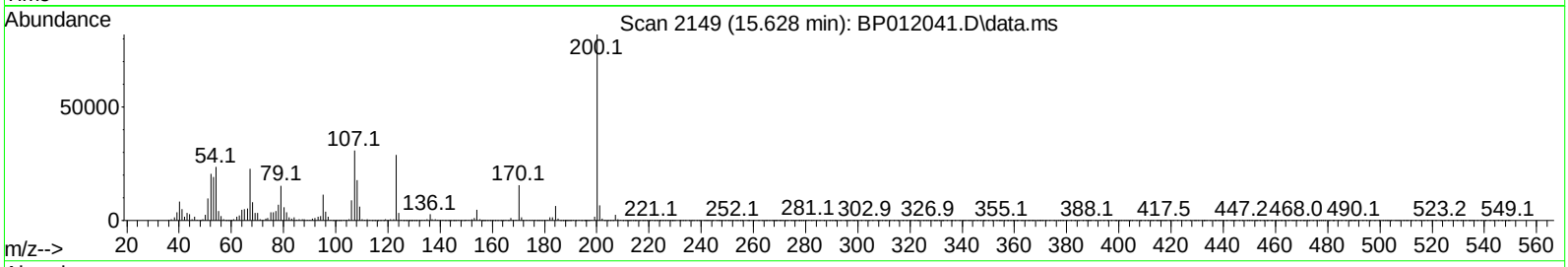

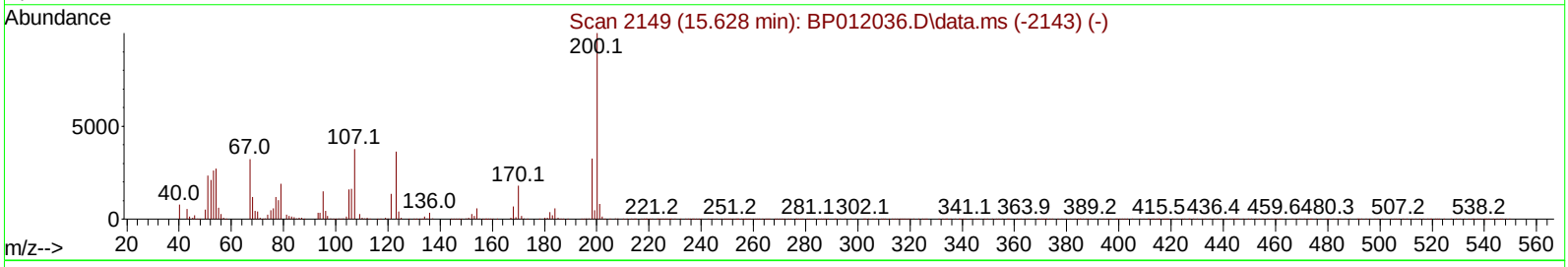

TIC: BP012041.D\data.ms

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

15.628min (-0.000) 20.87 ng/ul m

response 113972

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	17.90	18.99
52.00	34.50	25.15#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	223924	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	916610	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164	520234	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	1034392	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	975320	20.000	ng/ul	0.00
88) Perylene-d12	23.774	264	1051640	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	29815	5.250	ng/uL	0.00
4) Pyridine-d5	3.705	84	351823	23.293	ng/ul	0.00
7) Phenol-d5	7.028	99	480927	26.704	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	286076	27.985	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	394685	27.609	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	368714	26.655	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	183628	28.366	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	190015	28.948	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	333995	26.255	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	523023	27.899	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	970897	29.468	ng/ul	0.00
49) Acenaphthylene-d8	14.198	160	1143657	28.071	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	146962	22.720	ng/ul	0.00
60) Fluorene-d10	15.504	176	871435	29.234	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	113972m	20.867	ng/ul	0.00
73) Anthracene-d10	17.363	188	1335858	29.740	ng/ul	0.00
81) Pyrene-d10	19.598	212	1496051	29.941	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.616	264	1508079	29.499	ng/ul	0.00

Target Compounds	Qvalue

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

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