

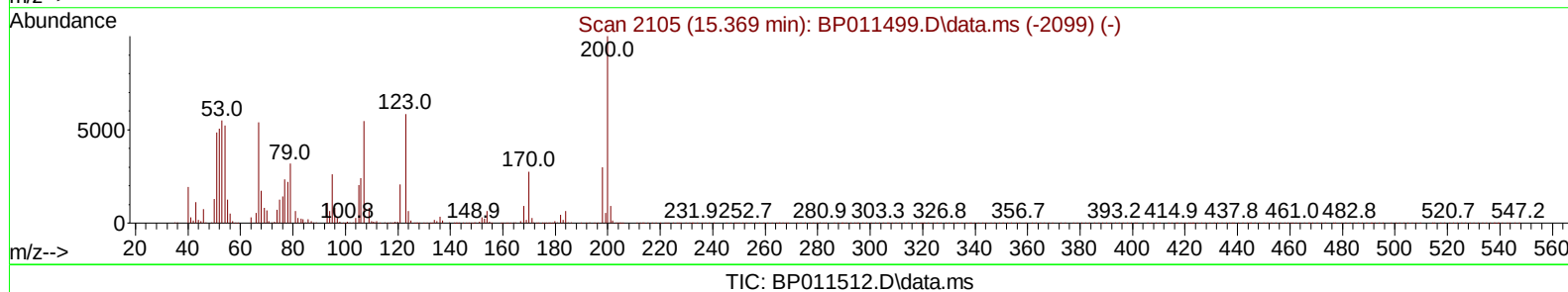
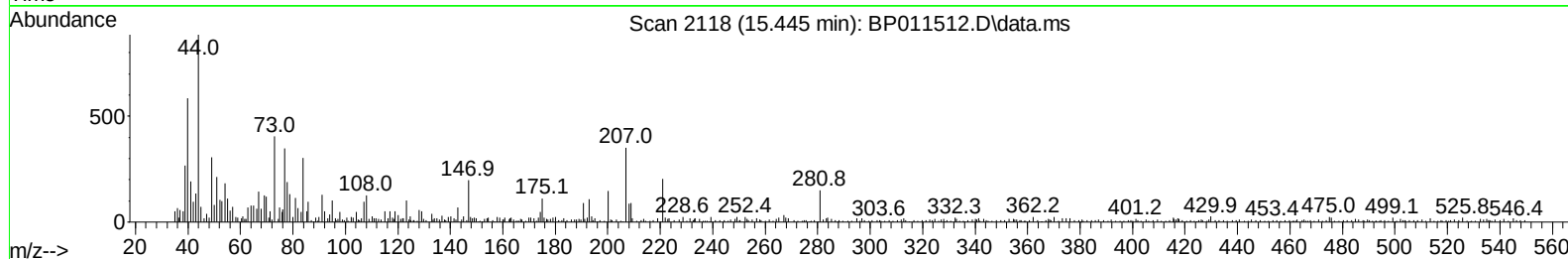
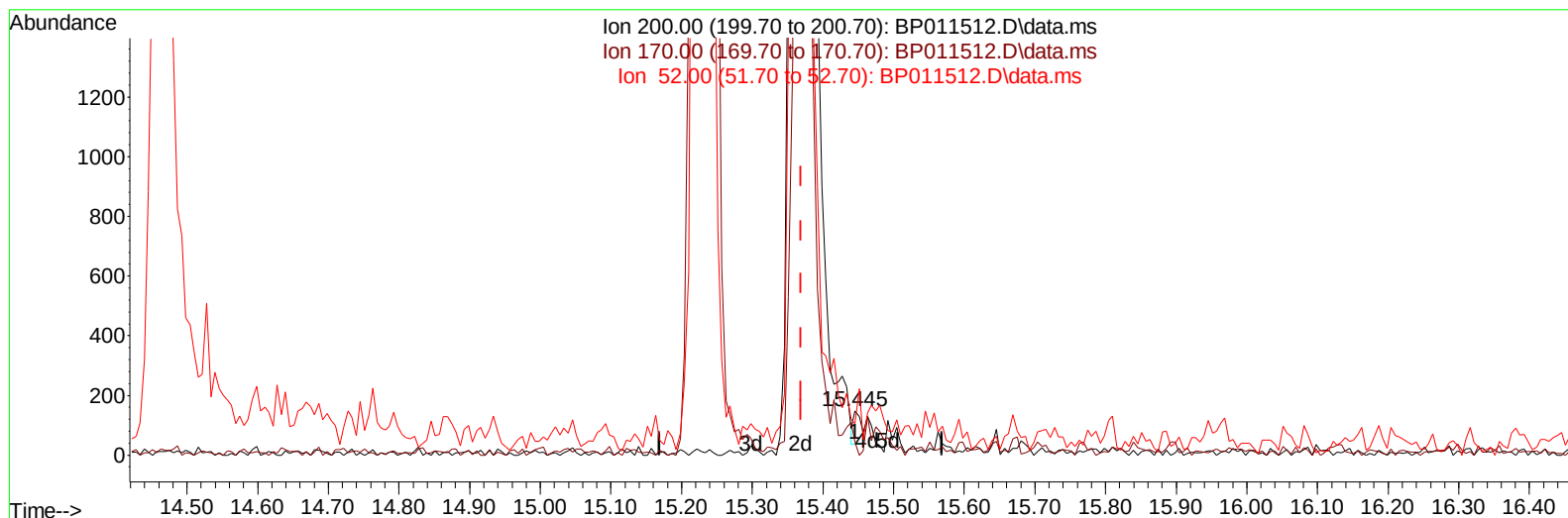
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011512.D
 Acq On : 19 Aug 2022 17:47
 Operator : CG/JU
 Sample : N4221-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KP9

Manual IntegrationsAPPROVED

Quant Time: Aug 20 04:38:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/22/2022
 Supervised By :Sohil Jodhani 08/22/2022



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

15.445min (+ 0.076) 0.04 ng/ul

response	84	
Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	29.25
52.00	67.40	71.43
0.00	0.00	0.00

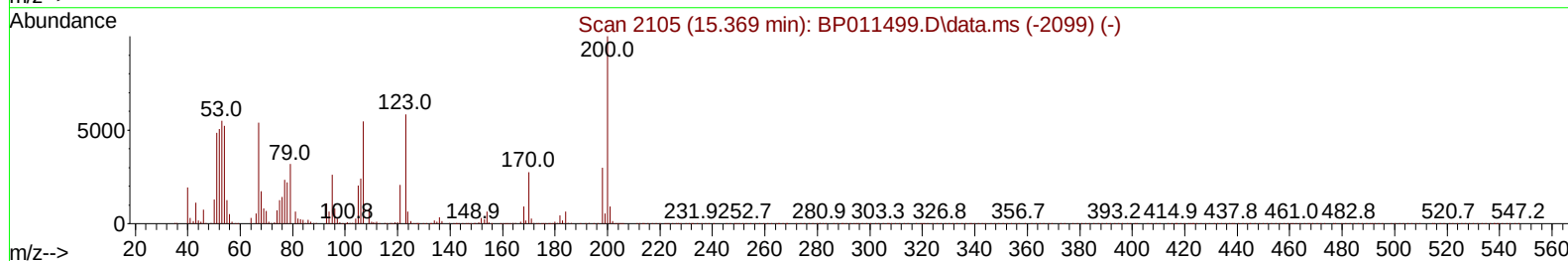
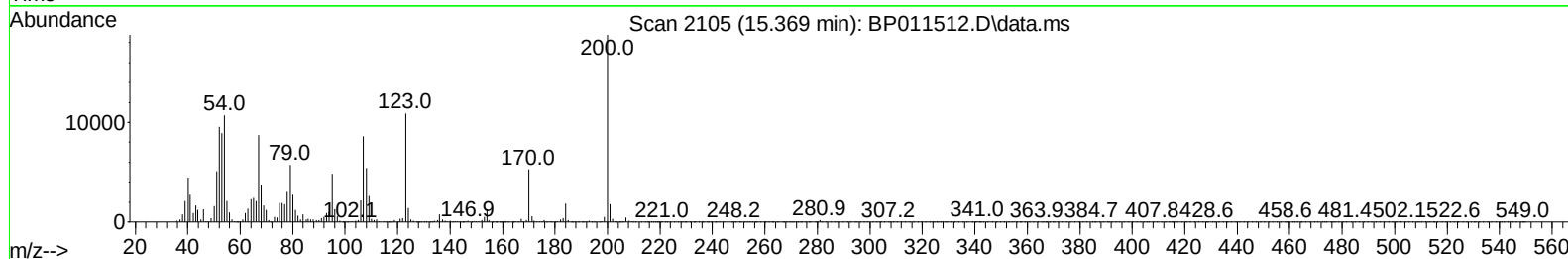
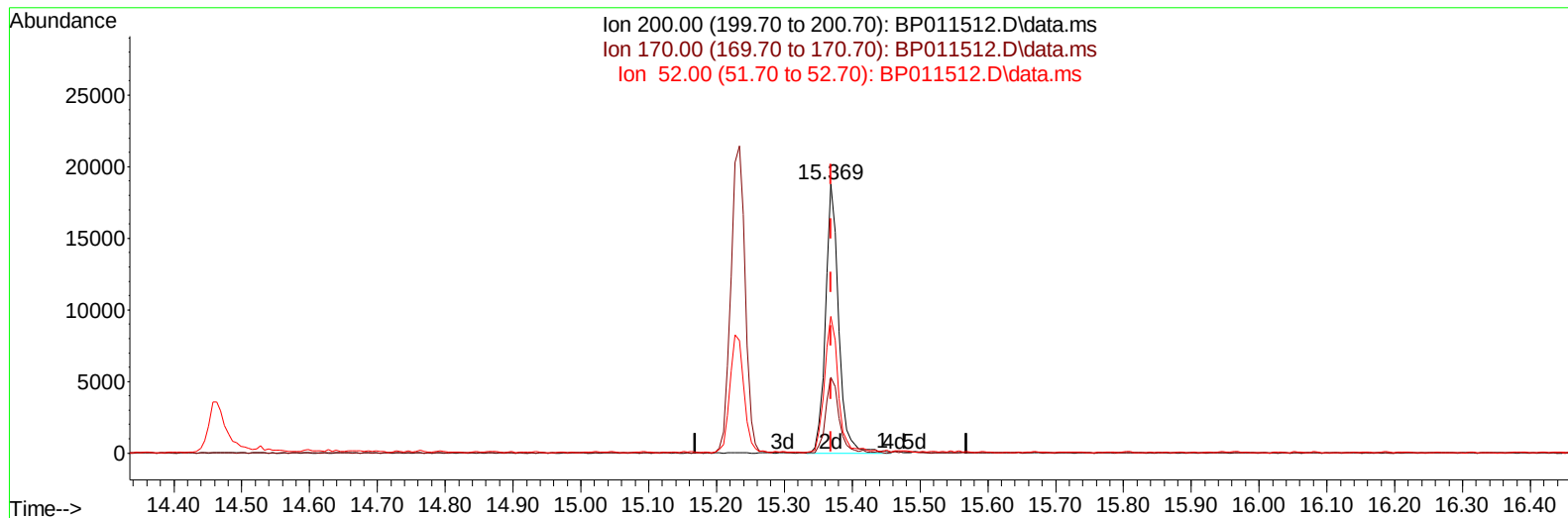
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TIC: BP011512.D\data.ms

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

15.369min (-0.000) 12.49 ng/ul m

response 25302

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	28.08
52.00	67.40	50.83#
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	7.557	152	86245	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	349456	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.228	164	188383	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	375022	20.000	ng/ul	0.00
79) Chrysene-d12	21.127	240	359840	20.000	ng/ul	0.00
88) Perylene-d12	23.421	264	375425	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	9224	3.319	ng/uL	0.00
4) Pyridine-d5	3.487	84	51513	7.239	ng/ul	0.00
7) Phenol-d5	6.740	99	145422	18.349	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	97360	18.908	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	120977	20.754	ng/ul	0.00
15) 4-Methylphenol-d8	8.281	113	115262	18.170	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	54778	21.773	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.434	143	55783	21.787	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	95360	20.177	ng/ul	0.00
31) 4-Chloroaniline-d4	10.498	131	140173	17.817	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	294196	22.080	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	329243	22.233	ng/ul	0.00
54) 4-Nitrophenol-d4	14.463	143	36540	14.234	ng/ul	0.00
60) Fluorene-d10	15.233	176	244081	21.775	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	25302m	12.486	ng/ul	0.00
73) Anthracene-d10	17.098	188	378487	22.788	ng/ul	0.00
81) Pyrene-d10	19.363	212	427718	23.493	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.268	264	422469	23.134	ng/ul	-0.01

Target Compounds Qvalue

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

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