

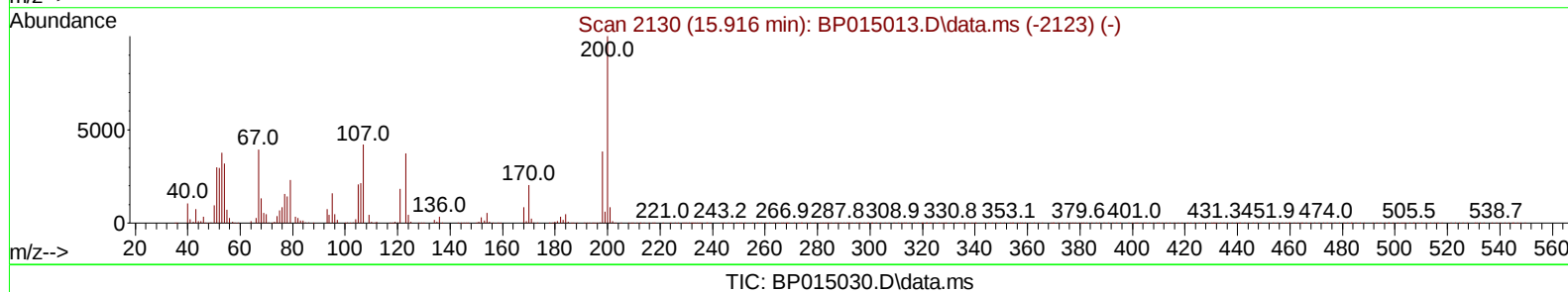
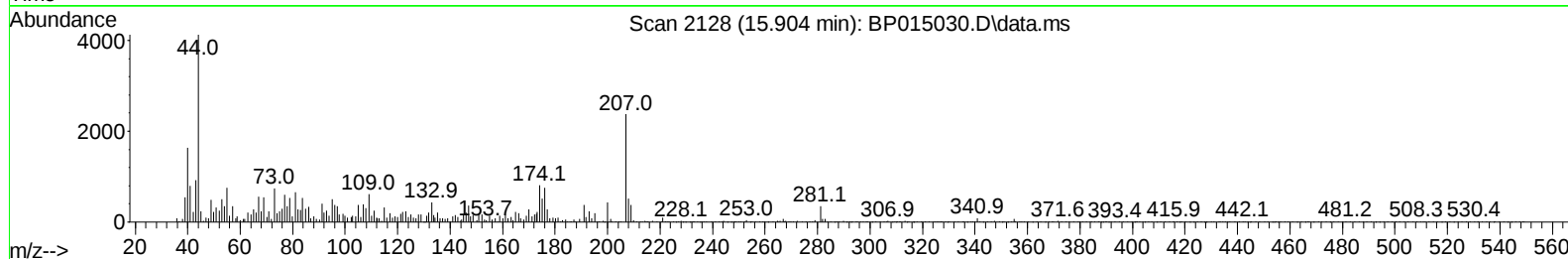
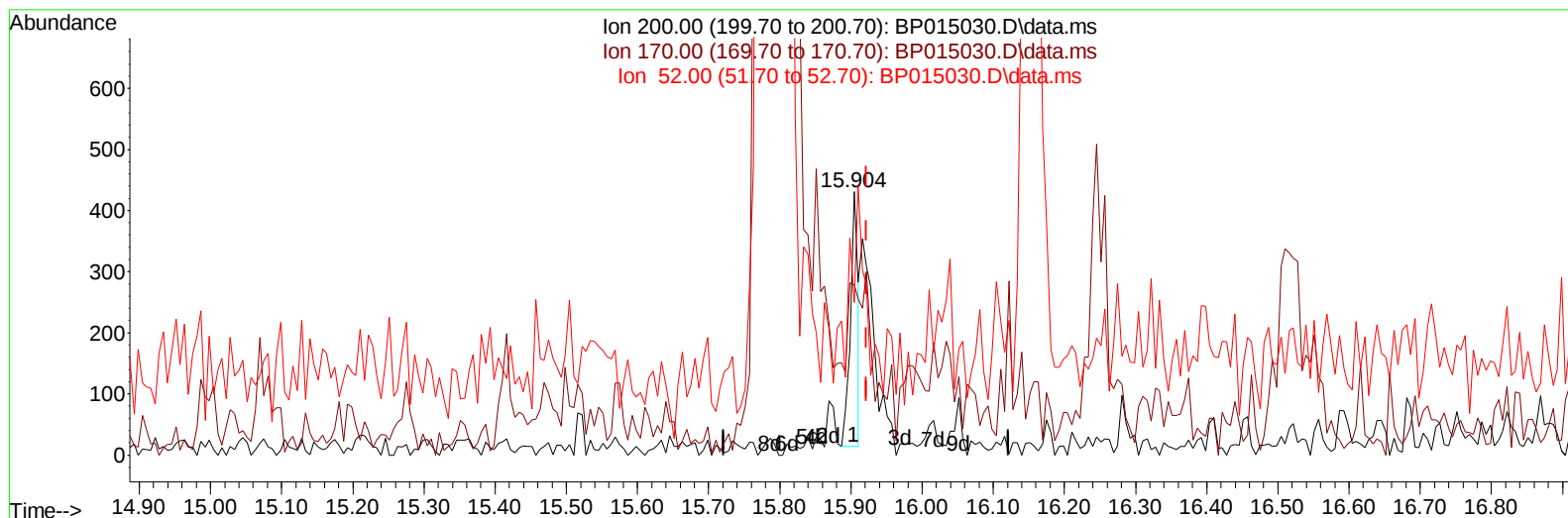
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050923\
Data File : BP015030.D
Acq On : 09 May 2023 22:01
Operator : CG/JU
Sample : 02609-24
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREL3

Manual IntegrationsAPPROVED

Quant Time: May 09 23:10:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 05 12:12:36 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/10/2023
Supervised By :Jagrut Upadhyay 05/10/2023



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

15.904min (-0.017) 0.02 ng/u1

response 319

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.30	64.73#
52.00	44.00	58.00#
0.00	0.00	0.00

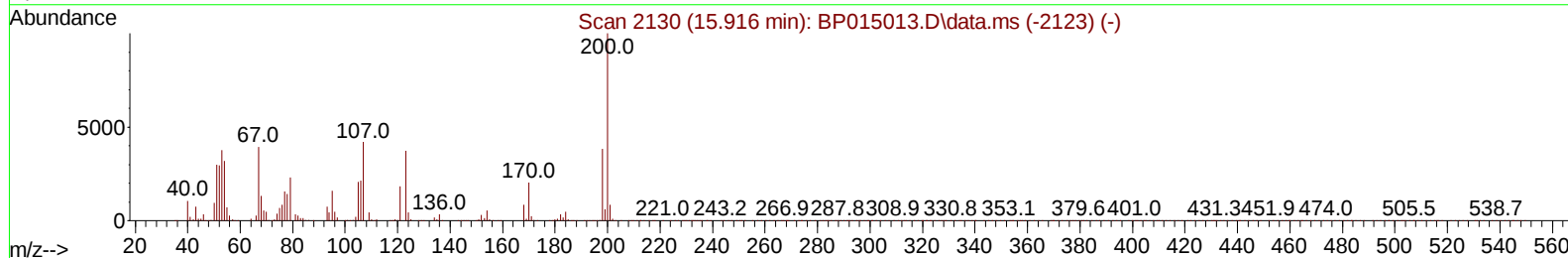
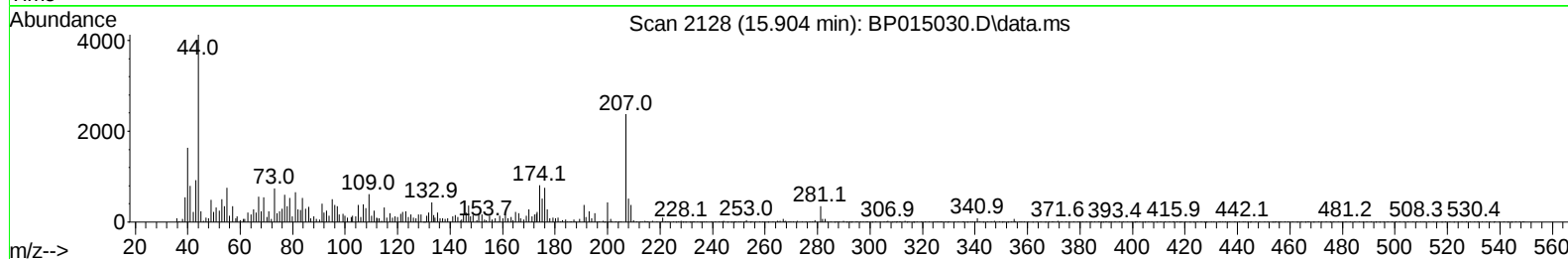
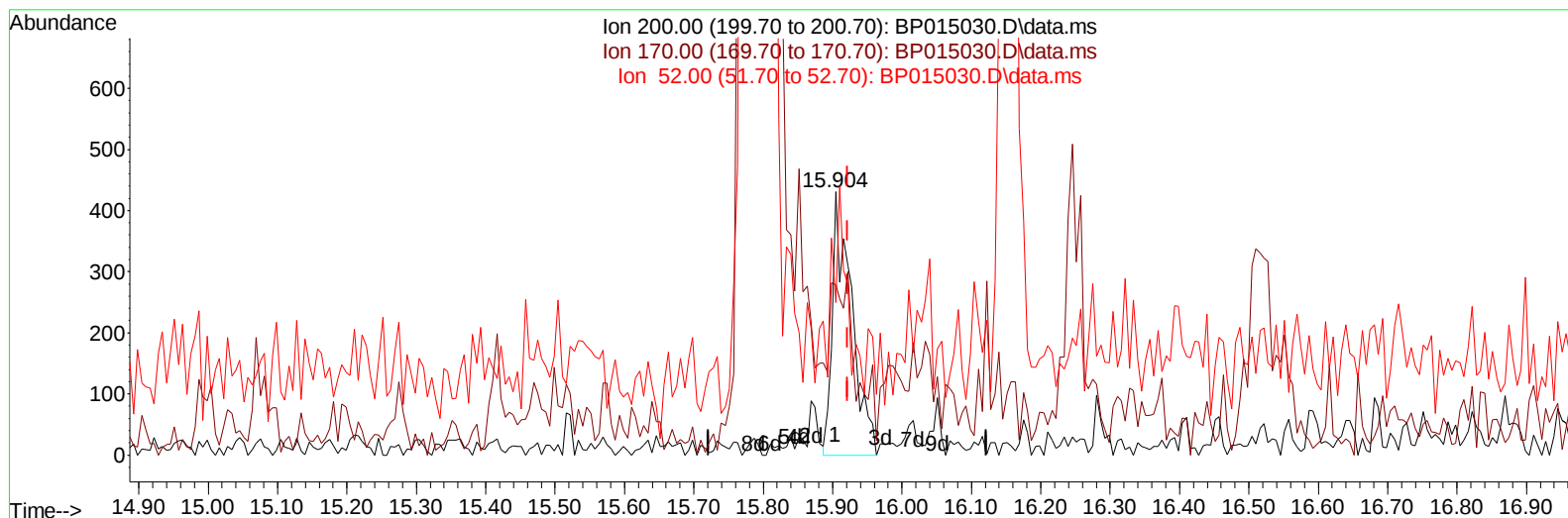
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TIC: BP015030.D\data.ms

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

15.904min (-0.017) 0.06 ng/ul m

response 828

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.30	64.73#
52.00	44.00	58.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050923\
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/10/2023
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Quant Time: May 09 23:11:51 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.158	152	393584	20.000	ng/ul	0.00
20) Naphthalene-d8	10.993	136	1882008	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.804	164	1239024	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.557	188	2601598	20.000	ng/ul	-0.01
79) Chrysene-d12	21.633	240	1744915	20.000	ng/ul	-0.01
88) Perylene-d12	24.227	264	1554535	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.458	96	35510	3.355	ng/uL	0.00
4) Pyridine-d5	3.899	84	115195	3.981	ng/ul	0.00
7) Phenol-d5	7.305	99	39567	1.160	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.475	67	490697	23.266	ng/ul	0.00
11) 2-Chlorophenol-d4	7.681	132	55120	2.197	ng/ul	0.00
15) 4-Methylphenol-d8	8.863	113	178803	6.800	ng/ul	-0.01
21) Nitrobenzene-d5	9.334	128	339807	25.387	ng/ul	0.00
24) 2-Nitrophenol-d4	10.063	143	14390	1.047	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.605	165	28794	1.067	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.128	131	332928	8.529	ng/ul	0.00
46) Dimethylphthalate-d6	14.204	166	2064058	23.538	ng/ul	0.00
49) Acenaphthylene-d8	14.498	160	2635738	25.062	ng/ul	0.00
54) 4-Nitrophenol-d4	14.987	143	206	0.014	ng/ul	-0.01
60) Fluorene-d10	15.793	176	1976048	26.192	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.904	200	828m	0.057	ng/ul	-0.02
73) Anthracene-d10	17.657	188	3096203	26.487	ng/ul	0.00
81) Pyrene-d10	19.875	212	3296651	28.012	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.057	264	2081755	27.484	ng/ul	-0.02

Target Compounds Qvalue

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

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