

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009871.D
 Acq On : 15 Apr 2022 21:14
 Operator : CG/JU
 Sample : PB144125BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

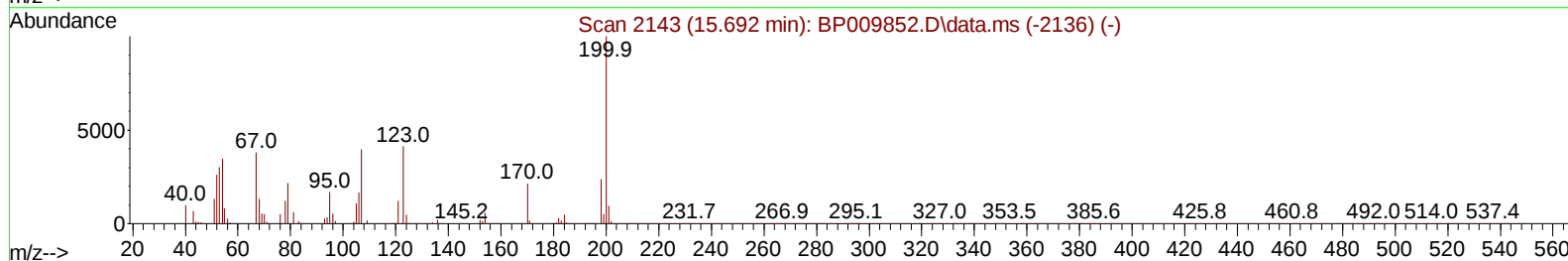
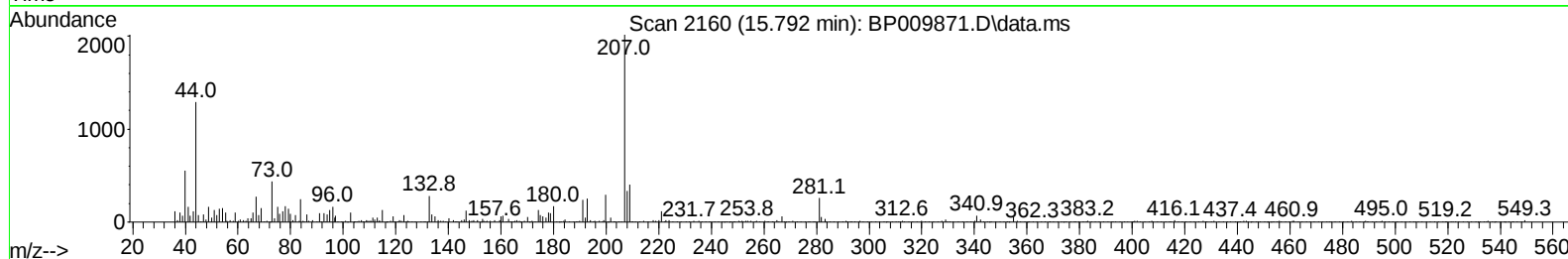
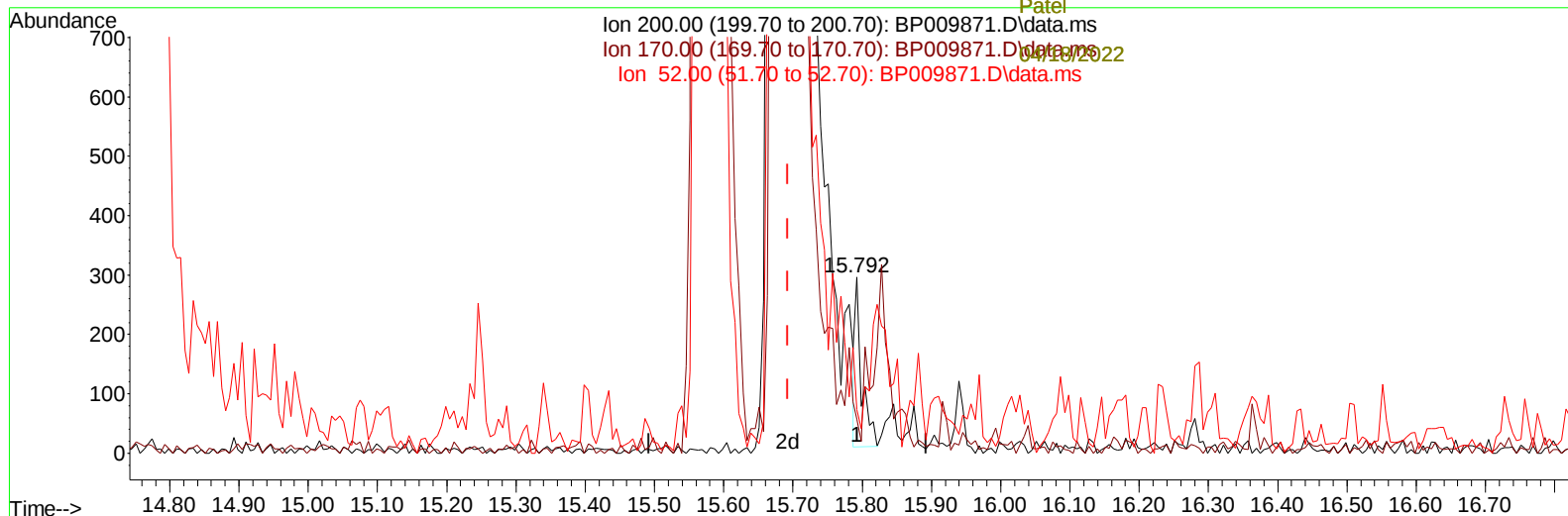
Instrument :
 BNA_P
 ClientSampleId :
 SBLK125

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:04:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

04/18/2022
 Supervised By :Yogesh
 Patel



TIC: BP009871.D\data.ms

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

15.792min (+ 0.100) 0.03 ng/u1

response 189

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	19.46
52.00	21.20	24.83
0.00	0.00	0.00

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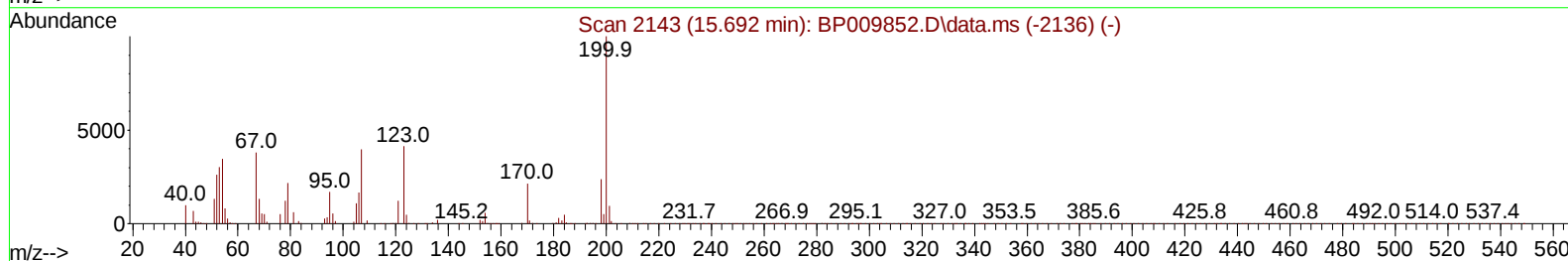
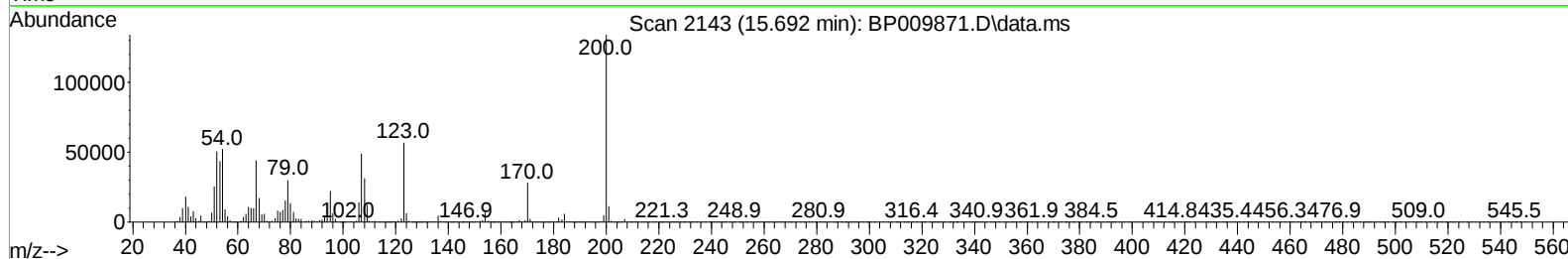
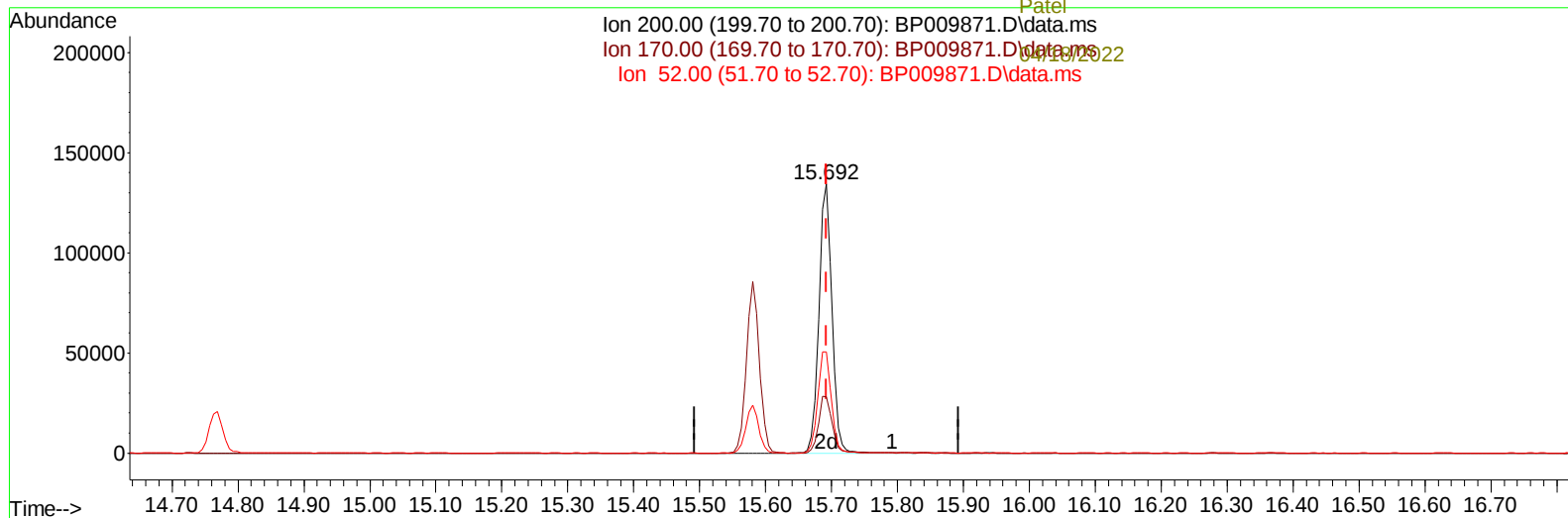
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.692min (+ 0.000) 32.51 ng/ul m

response 182845

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	20.95
52.00	21.20	37.60#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----04/18/2022							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.958	152		121773	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136		562139	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164		414407	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188		949537	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240		979446	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264		877061	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.381	96		18790	6.887	ng/uL	0.00
4) Pyridine-d5	3.799	84		266178	35.382	ng/ul	0.00
7) Phenol-d5	7.111	99		345874	32.352	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67		238671	37.499	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132		285926	35.456	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113		310521	34.794	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128		157454	38.842	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143		174639	39.295	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165		303397	32.623	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131		463439	34.987	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166		1204874	37.014	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160		1298648	33.674	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143		197123	33.391	ng/ul	0.00
60) Fluorene-d10	15.581	176		1004613	36.686	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200		182845m	32.505	ng/ul	0.00
73) Anthracene-d10	17.439	188		1690689	37.188	ng/ul	0.00
81) Pyrene-d10	19.669	212		2081136	38.905	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264		1829907	40.064	ng/ul	0.00

Target Compounds Qvalue

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

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