

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019004.D
 Acq On : 21 Dec 2023 17:57
 Operator : MA/JU
 Sample : PB157918BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

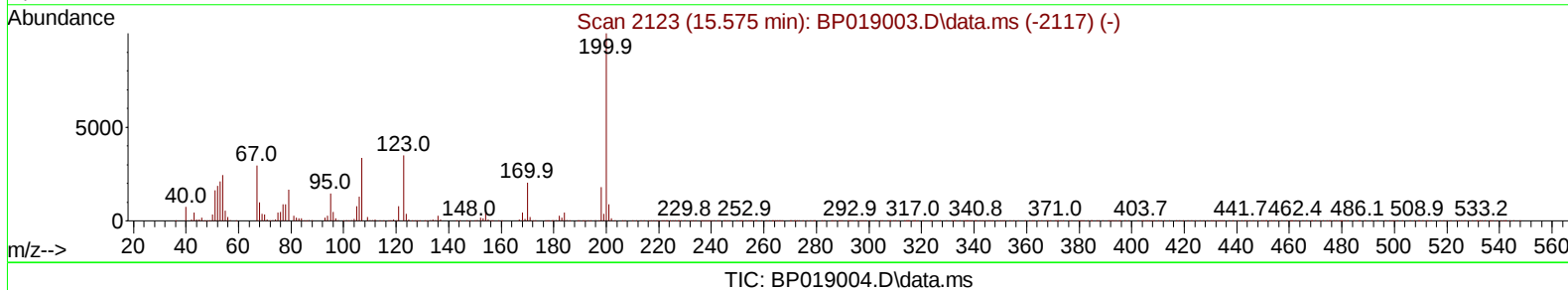
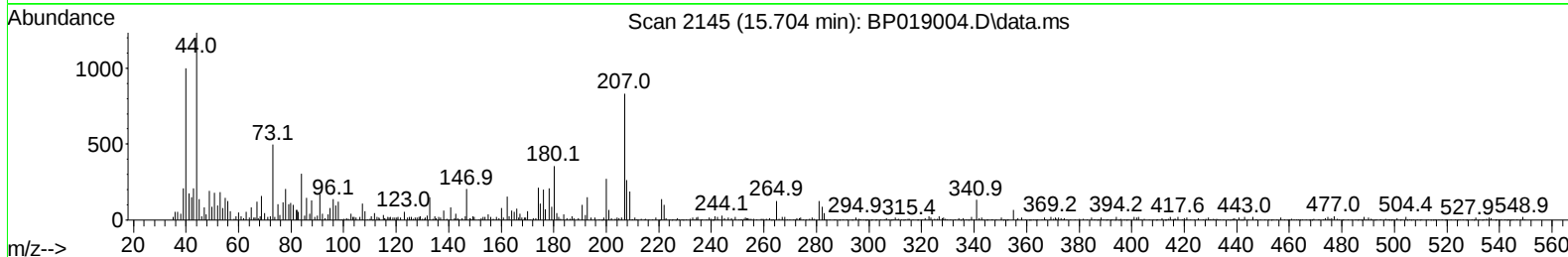
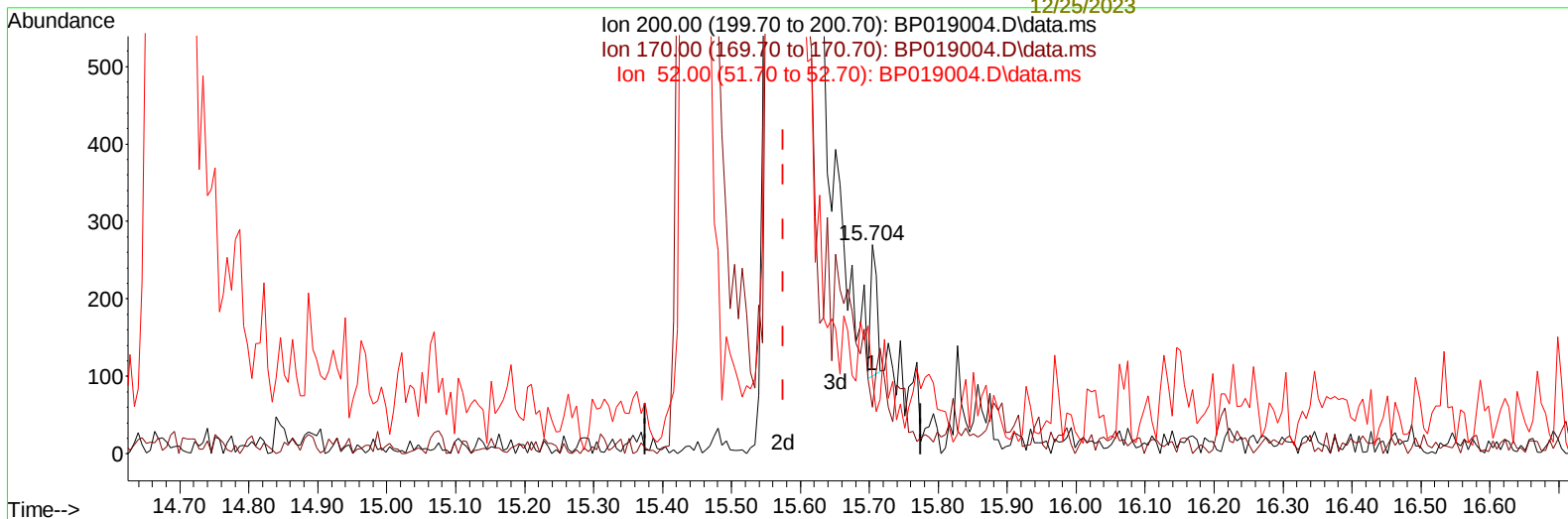
Instrument :
 BNA_P
 ClientSampleId :
 SBLK918

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/22/2023
 Supervised By :mohammad ahmed 12/25/2023

Supervised By :mohammad
 ahmed
 12/25/2023

Quant Time: Dec 22 08:47:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration



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

15.704min (+ 0.129) 0.03 ng/u1

response 107

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.50	22.22
52.00	36.60	35.56
0.00	0.00	0.00

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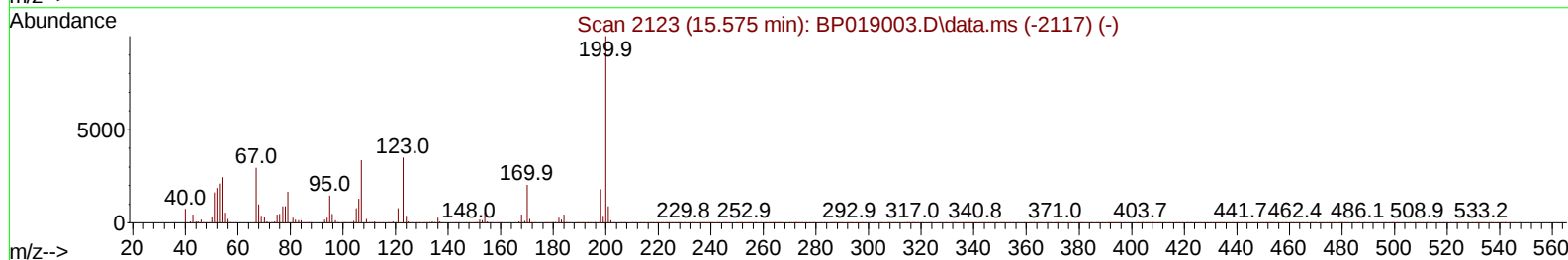
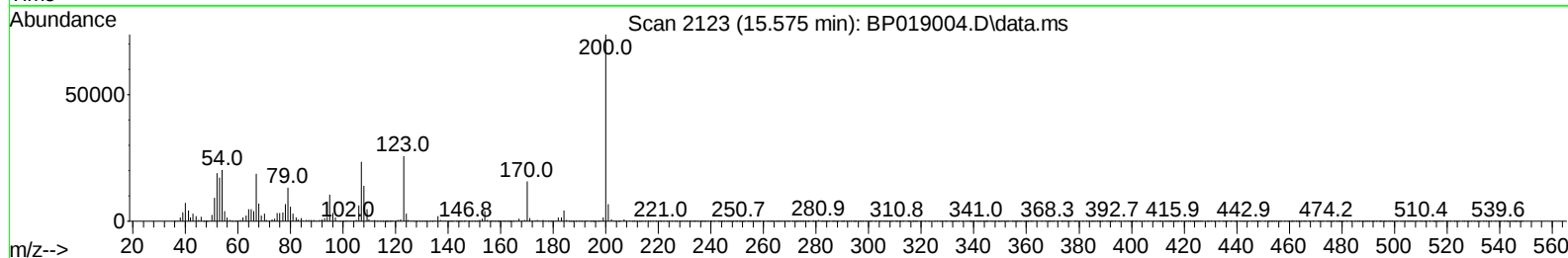
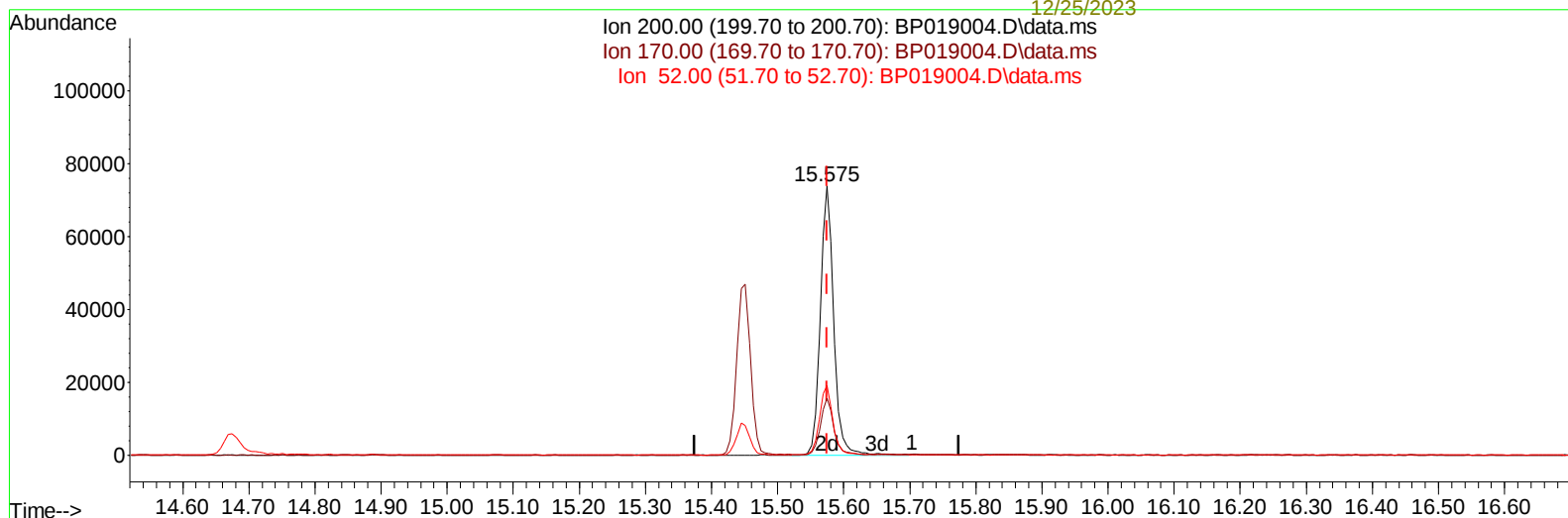
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TIC: BP019004.D\data.ms

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

15.575min (+ 0.000) 24.67 ng/ul m

response 104711

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.50	21.20
52.00	36.60	25.89#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.822	152		133191	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136		481387	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164		298619	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188		707936	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.298	240		823483	20.000	ng/ul	# 0.00
88) Perylene-d12	23.657	264		998054	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.252	96		18466	4.725	ng/uL	0.00
4) Pyridine-d5	3.669	84		239837	22.909	ng/ul	0.00
7) Phenol-d5	6.993	99		278134	20.809	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67		163203	20.601	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132		226711	21.737	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113		212404	19.636	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128		102046	23.817	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143		116991	26.337	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165		208272	22.930	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131		285820	24.078	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166		667166	25.033	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160		707374	24.103	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143		101480	23.196	ng/ul	0.00
60) Fluorene-d10	15.451	176		559689	25.018	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200		104711m	24.668	ng/ul	0.00
73) Anthracene-d10	17.304	188		941091	25.286	ng/ul	0.00
81) Pyrene-d10	19.545	212		1220770	19.044	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264		1499042	25.760	ng/ul	0.00

Target Compounds Qvalue

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

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