Quantitation Report (Qedit)

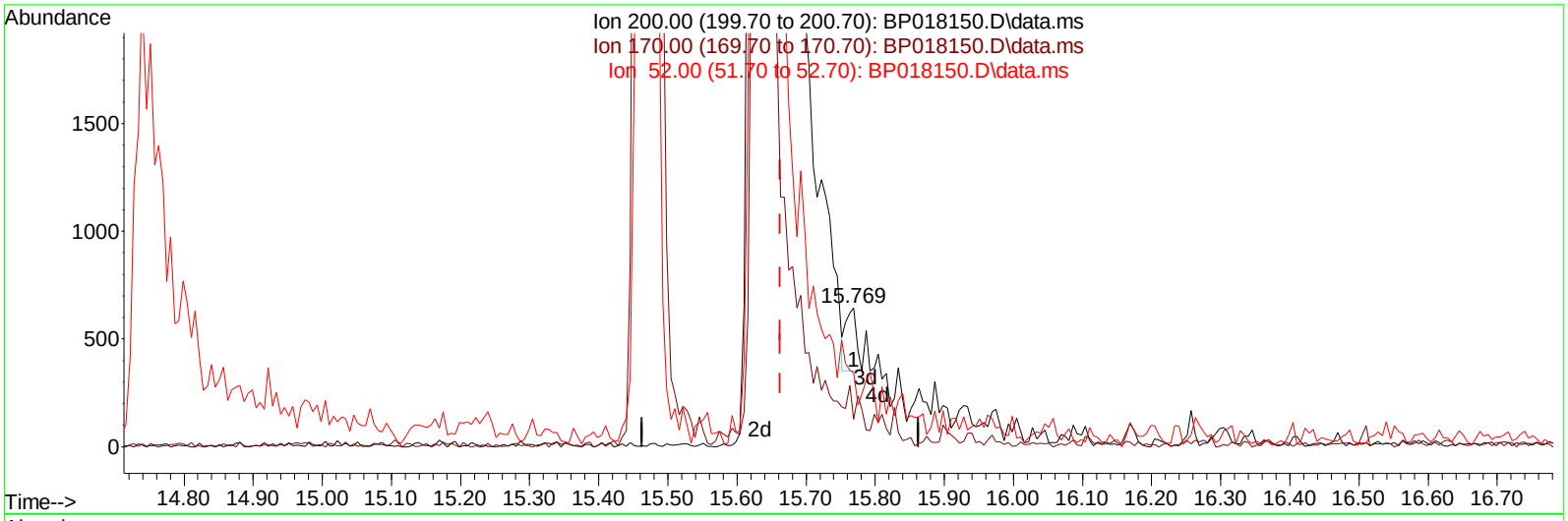
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018150.D
Acq On : 14 Nov 2023 11:39
Operator : MA/JU
Sample : PB156806BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

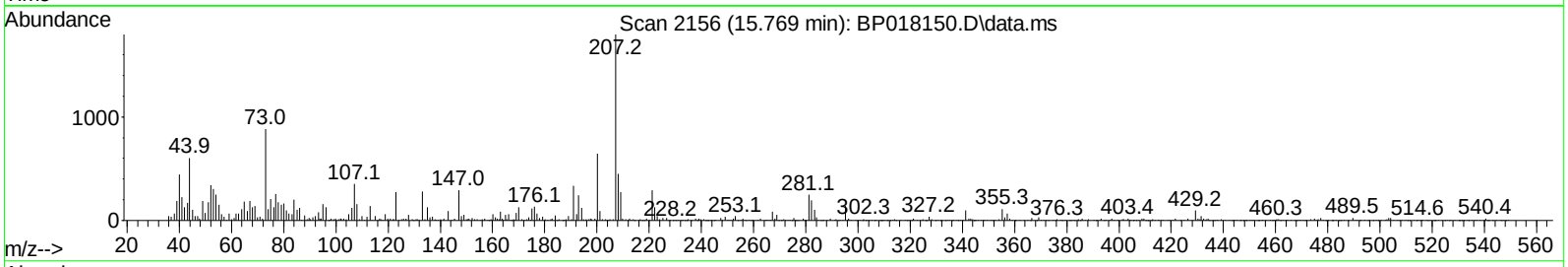
Quant Time: Nov 15 00:37:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

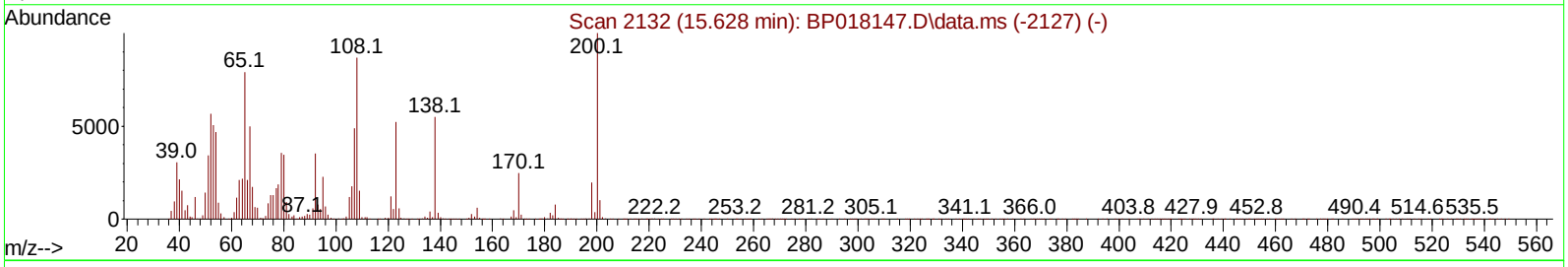
Instrument :
BNA_P
ClientSampleId :
SBLK806

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023







TIC: BP018150.D\data.ms

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

15.769min (+ 0.106) 0.11 ng/ul

response 312

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	19.57
52.00	52.00	53.26
0.00	0.00	0.00

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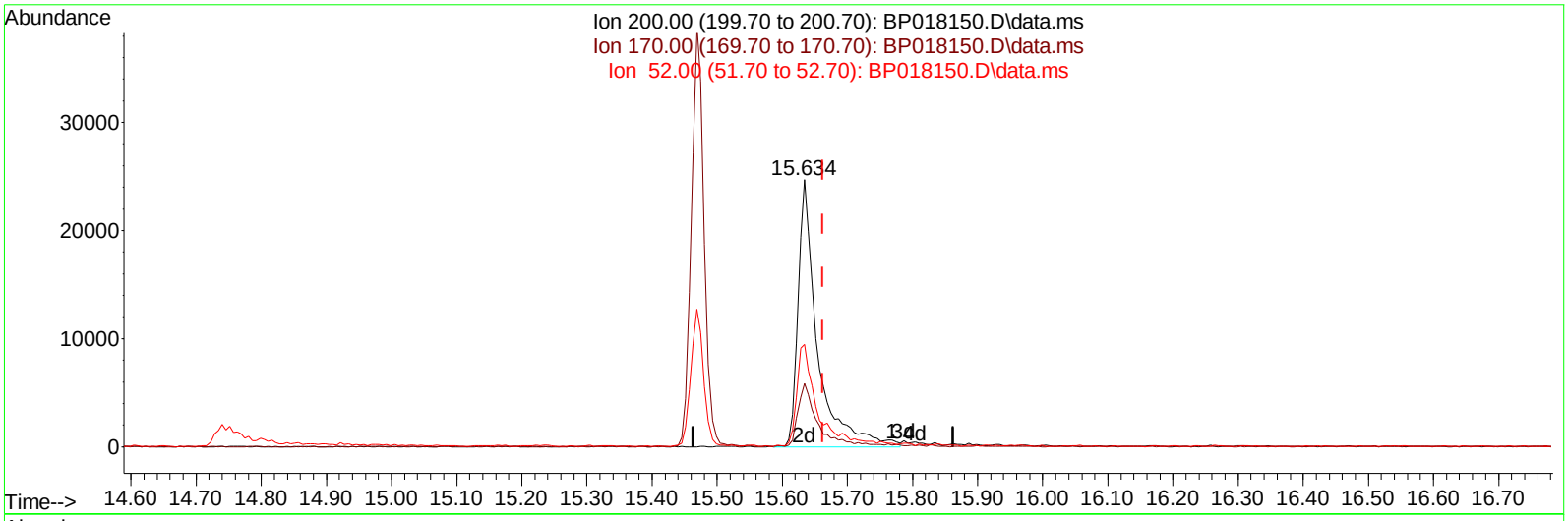

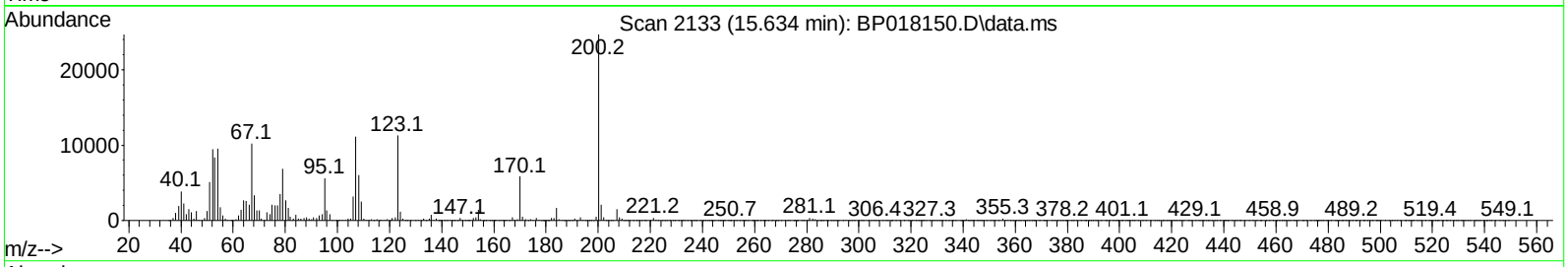

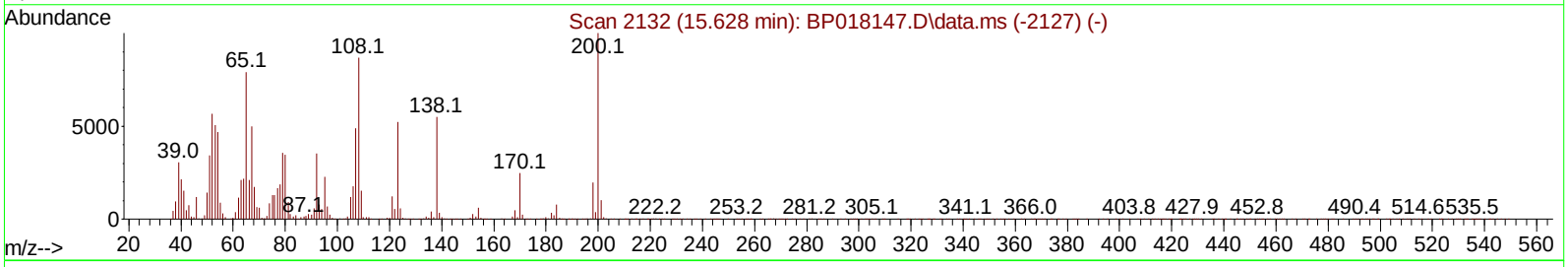

TIC: BP018150.D\data.ms

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

15.634min (-0.029) 18.71 ng/ul m

response 50853

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	23.66
52.00	52.00	38.27#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Quant Time: Nov 15 00:58:07 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	77789	20.000	ng/ul	0.00
20) Naphthalene-d8	10.605	136	322273	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.469	164	162799	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	375951	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	391066	20.000	ng/ul	#-0.01
88) Perylene-d12	23.845	264	406027	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	14808	6.715	ng/uL	0.00
4) Pyridine-d5	3.599	84	128539	22.356	ng/ul	0.00
7) Phenol-d5	6.958	99	228937	30.484	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	155822	35.307	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	189020	31.783	ng/ul	0.00
15) 4-Methylphenol-d8	8.511	113	186072	30.236	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	89570	30.849	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	98142	29.836	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	163970	27.968	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	214042	26.622	ng/ul	0.00
46) Dimethylphthalate-d6	13.893	166	523650	34.861	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	530238	32.409	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	37263	16.601	ng/ul	0.00
60) Fluorene-d10	15.469	176	418517	33.747	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	50853m	18.711	ng/ul	-0.03
73) Anthracene-d10	17.351	188	683869	33.634	ng/ul	0.00
81) Pyrene-d10	19.604	212	950967	36.405	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.680	264	826080	34.990	ng/ul	-0.02
Target Compounds						Qvalue

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

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