Quantitation Report (Qedit)

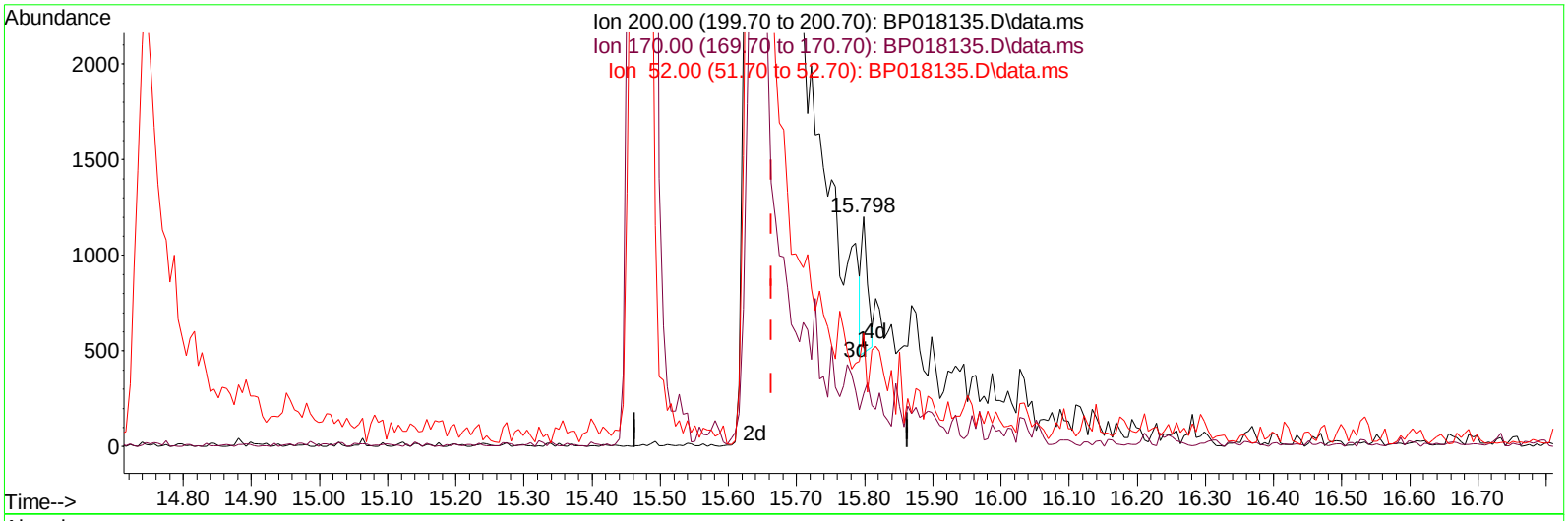
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111323\
Data File : BP018135.D
Acq On : 13 Nov 2023 17:25
Operator : MA/JU
Sample : PB157057BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

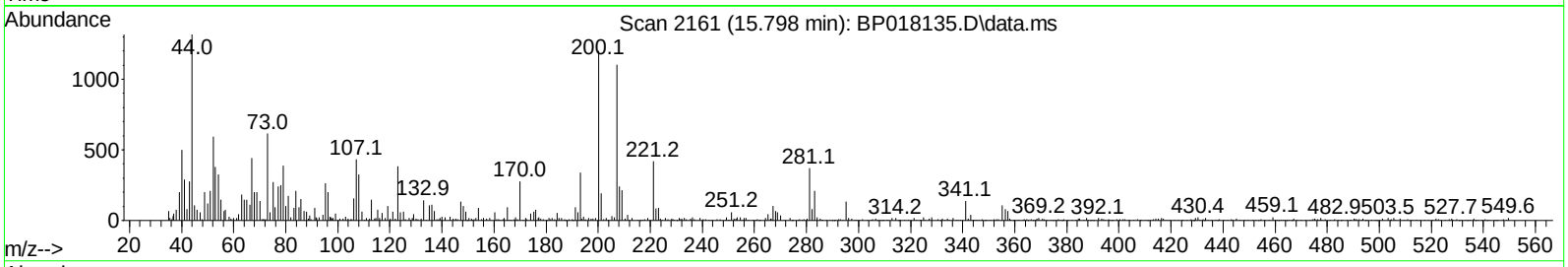
Quant Time: Nov 13 19:17:44 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

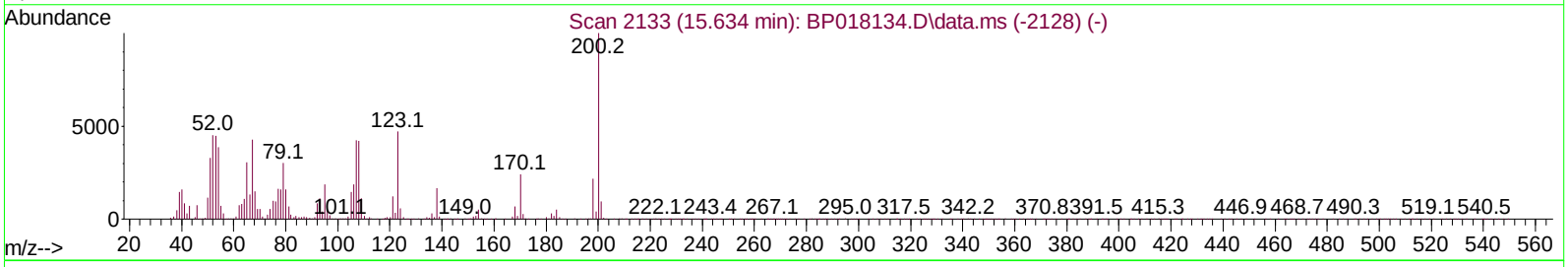
Instrument :
BNA_P
ClientSampleId :
SBLK057

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/14/2023
Supervised By :mohammad ahmed 11/16/2023







TIC: BP018135.D\data.ms

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

15.798min (+ 0.135) 0.12 ng/ul

response 417

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	23.06
52.00	52.00	49.54
0.00	0.00	0.00

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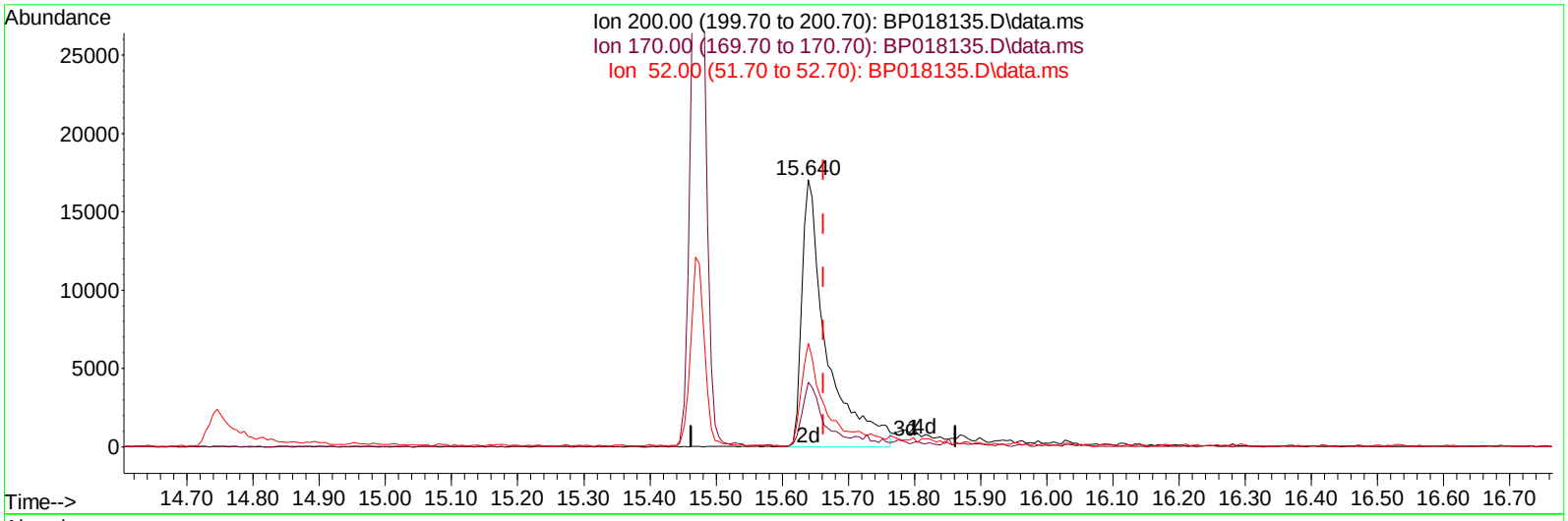

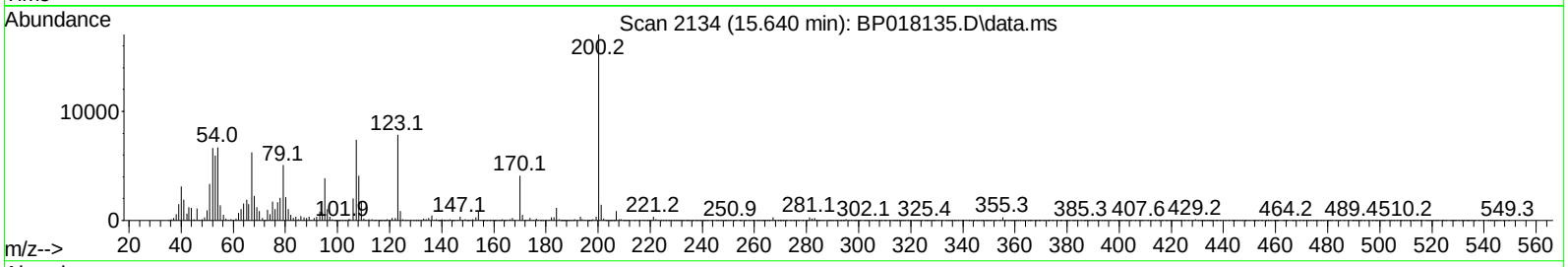

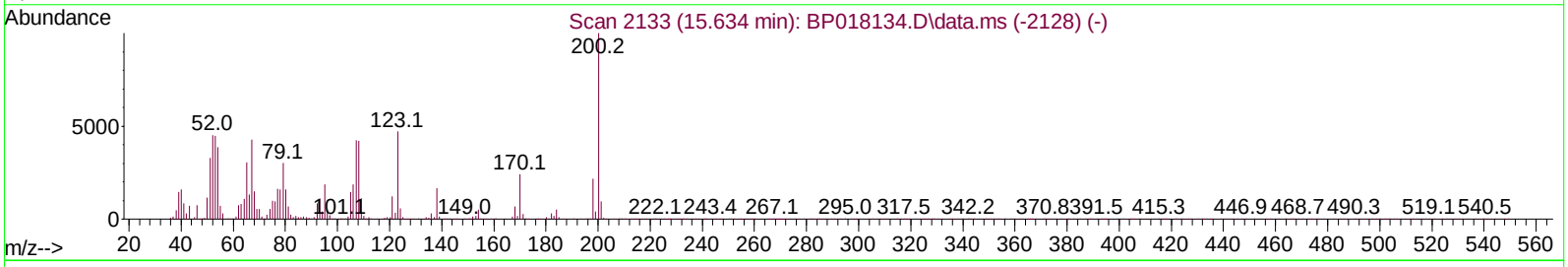

TIC: BP018135.D\data.ms

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

15.640min (-0.023) 12.99 ng/ul m

response 44203

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	24.22
52.00	52.00	38.83#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	82540	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	290312	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	210997	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	470557	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	423874	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	454410	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	13299	5.684	ng/uL	0.00
4) Pyridine-d5	3.605	84	99872	16.370	ng/ul	0.00
7) Phenol-d5	6.958	99	220436	27.663	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	149275	31.876	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	177160	28.074	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	170085	26.048	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	87570	33.480	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	82211	27.744	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	126663	23.983	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	156469	21.604	ng/ul	0.00
46) Dimethylphthalate-d6	13.893	166	605906	31.123	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	614832	28.996	ng/ul	0.00
54) 4-Nitrophenol-d4	14.746	143	37662	12.946	ng/ul	0.00
60) Fluorene-d10	15.475	176	482117	29.995	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	44203m	12.994	ng/ul	-0.02
73) Anthracene-d10	17.351	188	789180	31.010	ng/ul	0.00
81) Pyrene-d10	19.610	212	916859	32.383	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	810994	30.693	ng/ul	-0.02
Target Compounds						Qvalue

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

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