Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032923\
Data File : BP014384.D
Acq On : 30 Mar 2023 03:43
Operator : CG/JU
Sample : PB151679BL
Misc :
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Mar 30 04:23:01 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 30 00:56:20 2023
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
SBLK679

Manual IntegrationsAPPROVED

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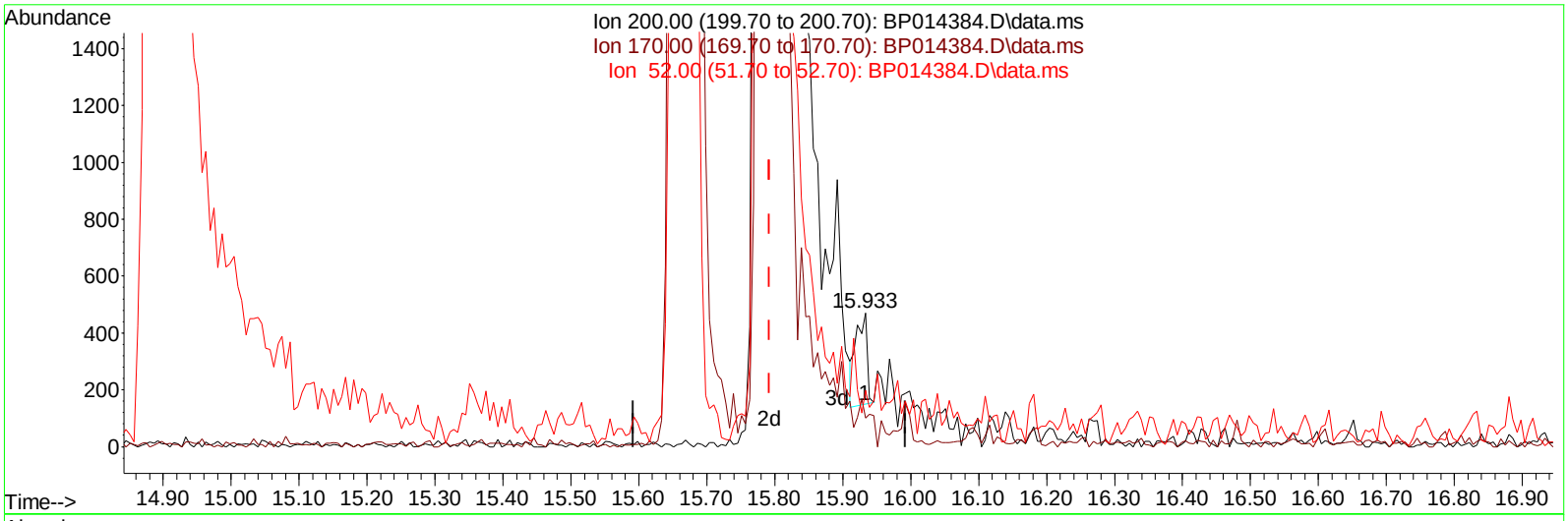
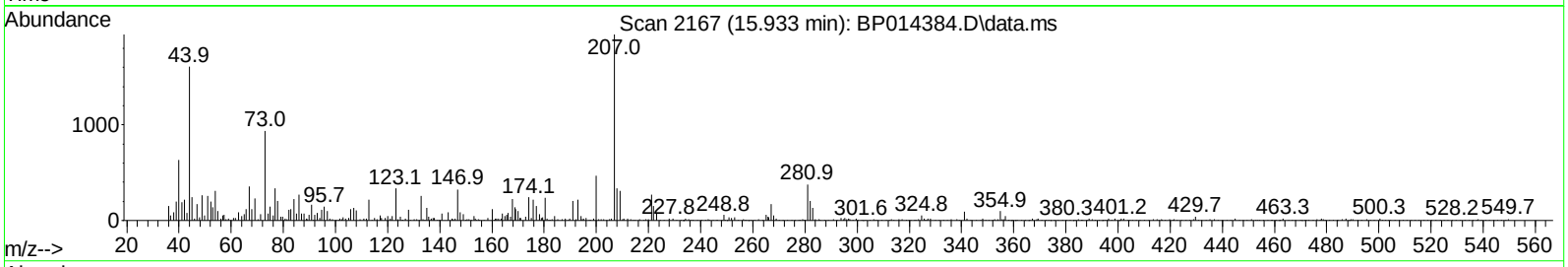
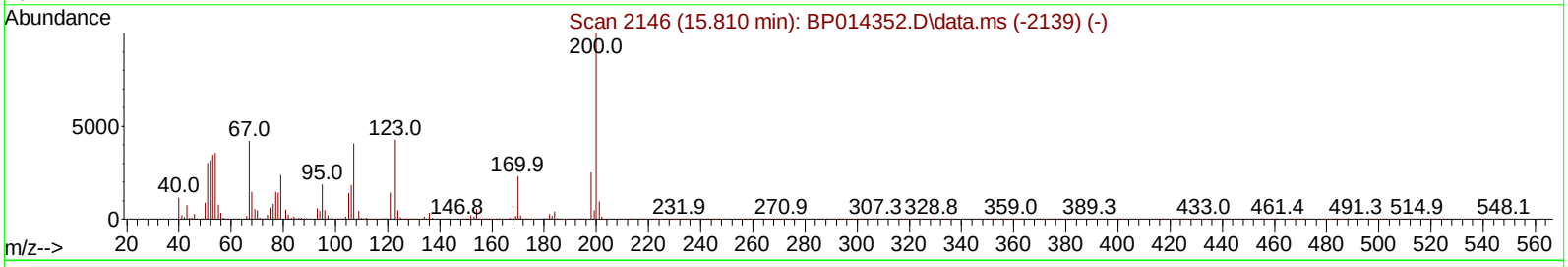
Supervised By :Jagrut Upadhyay 03/30/2023







TIC: BP014384.D\data.ms

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

15.933min (+ 0.141) 0.05 ng/ul

response 378

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	21.23
52.00	45.50	42.04
0.00	0.00	0.00

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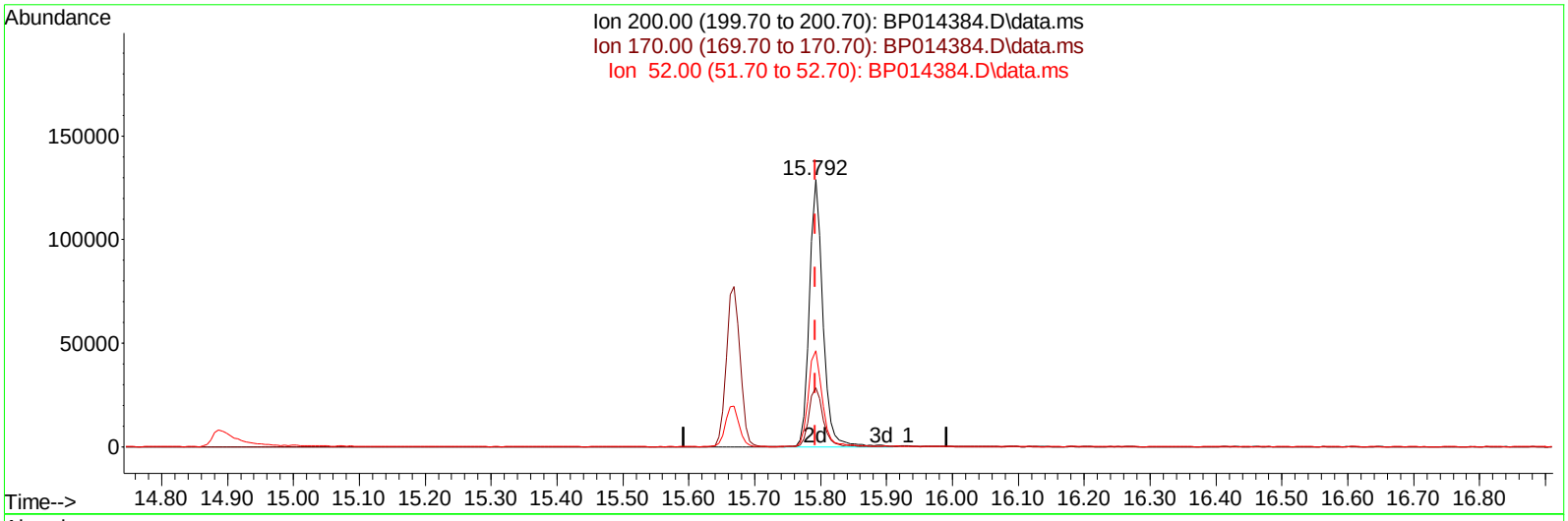

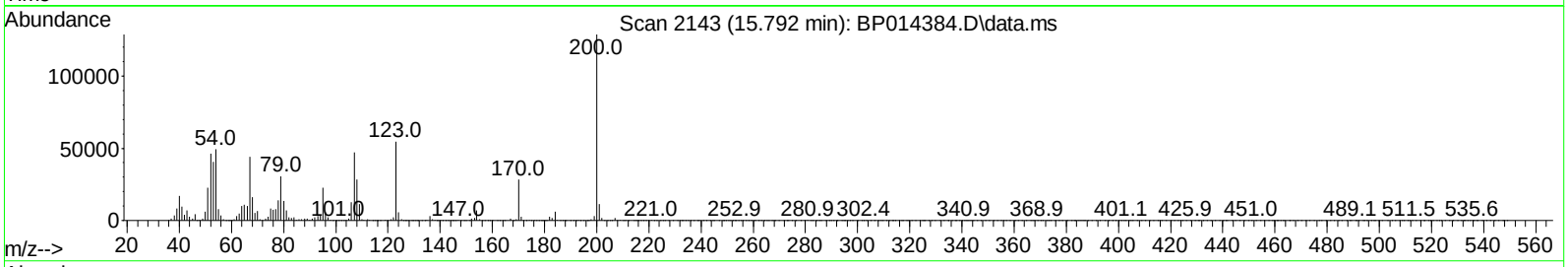

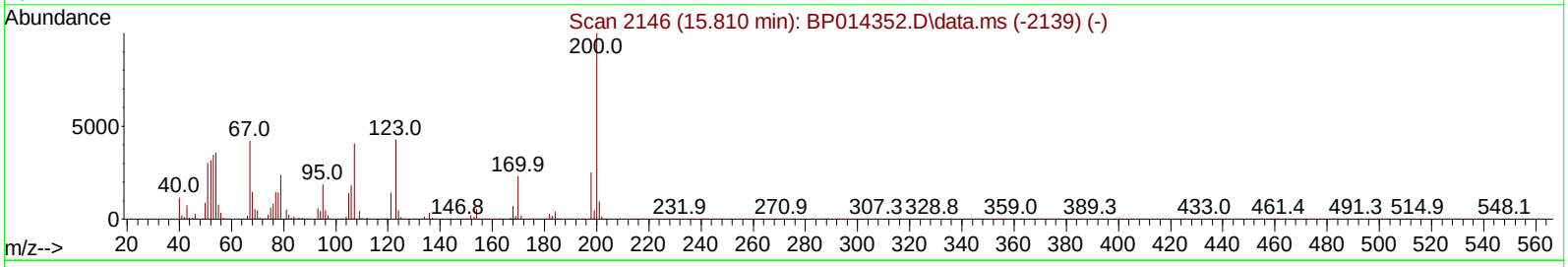

TIC: BP014384.D\data.ms

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

15.792min (+ 0.000) 25.08 ng/ul m

response 182190

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	22.02
52.00	45.50	35.98#
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	8.016	152	143130	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	614375	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	459621	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.433	188	987328	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	958827	20.000	ng/ul	0.00
88) Perylene-d12	24.045	264	917666	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	21702	5.880	ng/uL	0.00
4) Pyridine-d5	3.822	84	322366	34.082	ng/ul	0.00
7) Phenol-d5	7.181	99	388209	29.409	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	249386	29.446	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	283295	29.694	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	315602	29.641	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	143217	28.876	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	166854	29.379	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	287054	27.753	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	399952	29.357	ng/ul	0.00
46) Dimethylphthalate-d6	14.081	166	1072062	28.767	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	1160882	28.074	ng/ul	0.00
54) 4-Nitrophenol-d4	14.886	143	150425	26.718	ng/ul	0.00
60) Fluorene-d10	15.669	176	928103	29.864	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	182190m	25.078	ng/ul	0.00
73) Anthracene-d10	17.533	188	1449062	30.326	ng/ul	0.00
81) Pyrene-d10	19.763	212	1786527	27.739	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	1561538	32.028	ng/ul	0.00

Target Compounds Qvalue

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

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