Quantitation Report (Qedit)

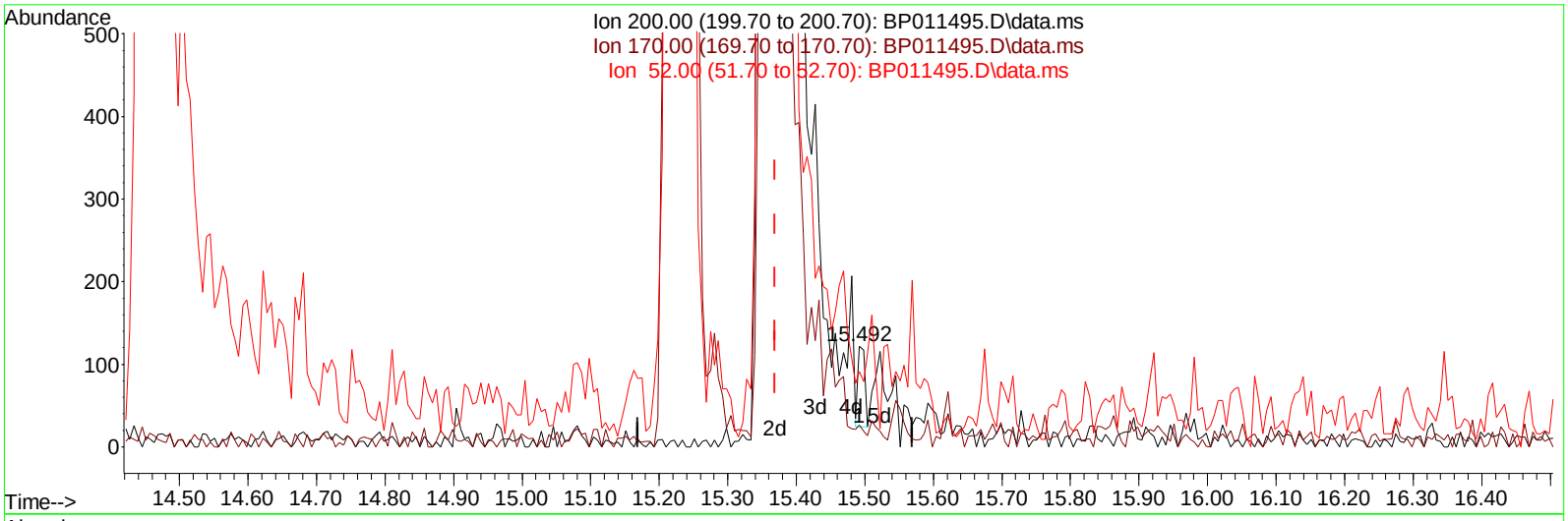
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
Data File : BP011495.D
Acq On : 18 Aug 2022 23:15
Operator : CG/JU
Sample : PB147098BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

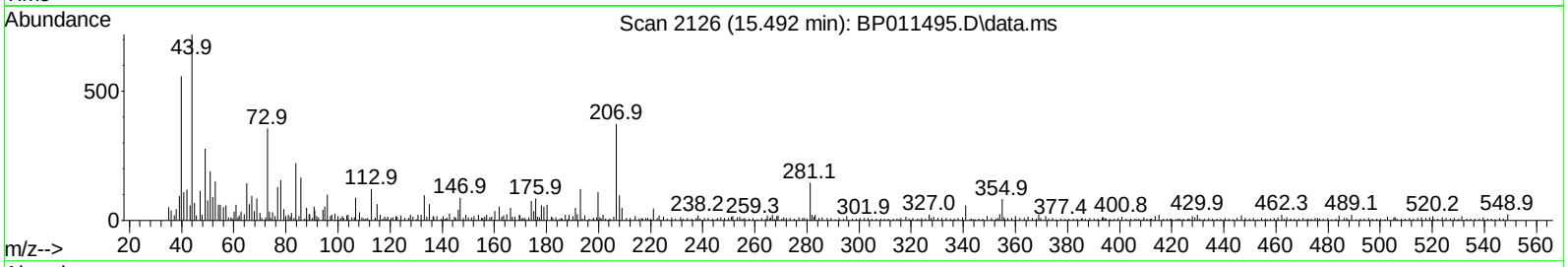
Quant Time: Aug 19 00:20:23 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 01:58:37 2022
Response via : Initial Calibration

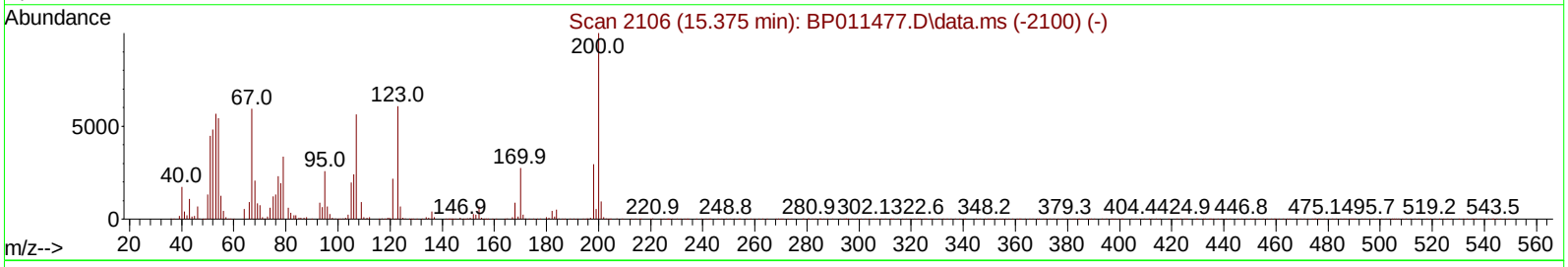
Instrument :
BNA_P
ClientSampleId :
SBLK098

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/19/2022
Supervised By :Sohil Jodhani 08/22/2022







TIC: BP011495.D\data.ms

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

15.492min (+ 0.123) 0.03 ng/ul

response 67

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	22.13
52.00	67.40	74.59
0.00	0.00	0.00

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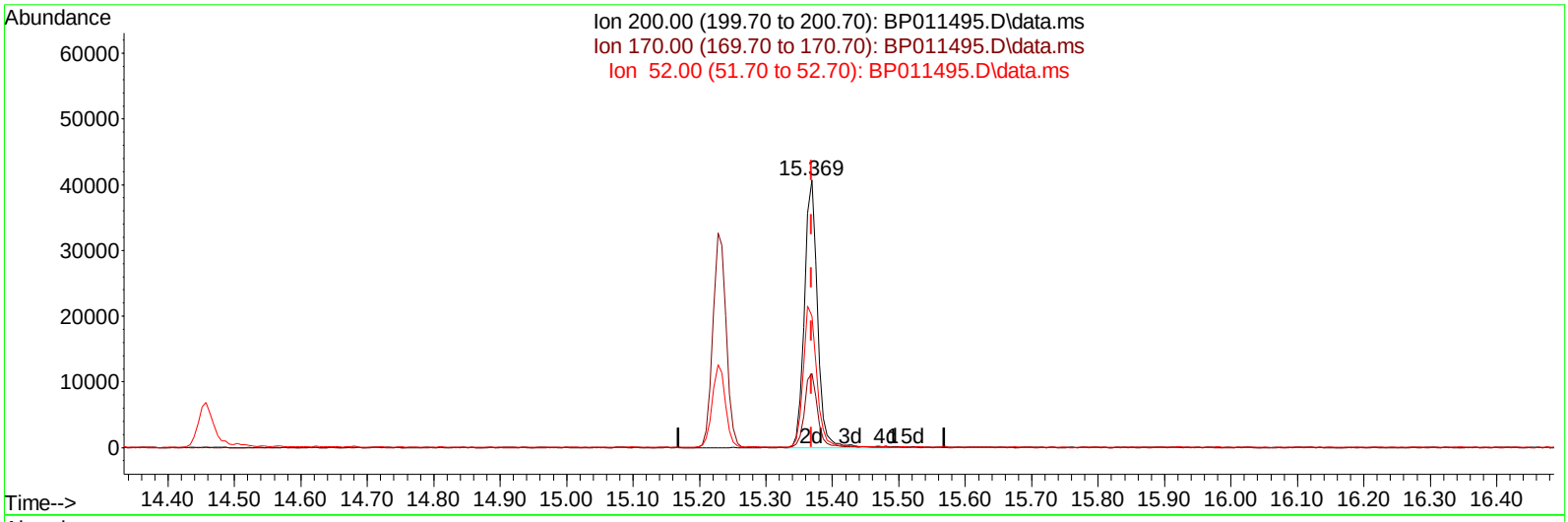

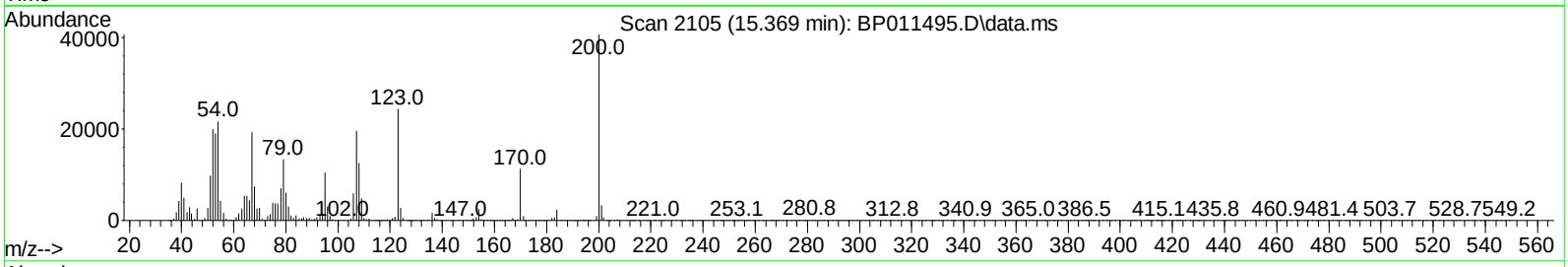

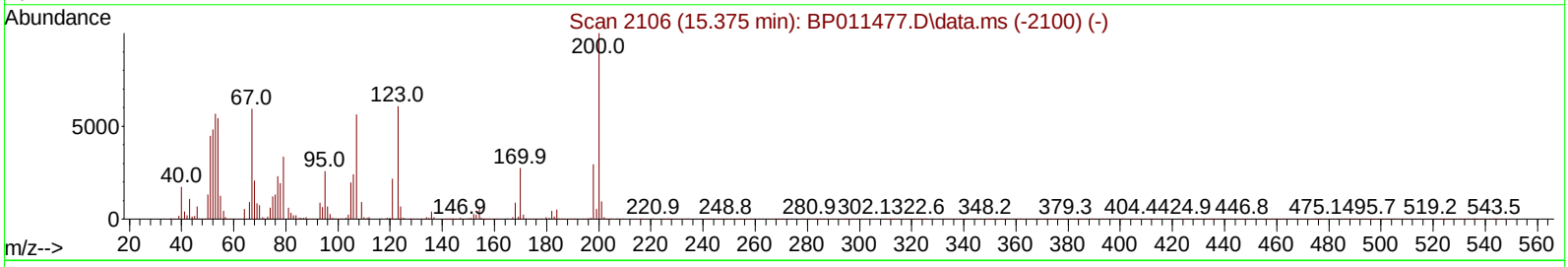

TIC: BP011495.D\data.ms

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

15.369min (-0.000) 23.67 ng/ul m

response 55122

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	27.71
52.00	67.40	49.09#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	100296	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	411423	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.228	164	225737	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	431063	20.000	ng/ul	0.00
79) Chrysene-d12	21.127	240	396494	20.000	ng/ul	0.00
88) Perylene-d12	23.421	264	435775	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	14977	4.634	ng/uL	0.00
4) Pyridine-d5	3.475	84	193571	23.393	ng/ul	0.00
7) Phenol-d5	6.740	99	215120	23.340	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.904	67	143108	23.899	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	176476	26.033	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	167632	22.723	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	81540	27.529	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.434	143	82175	27.260	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	138906	24.964	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	215570	23.273	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	433827	27.172	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	463923	26.144	ng/ul	0.00
54) 4-Nitrophenol-d4	14.457	143	63736	20.719	ng/ul	0.00
60) Fluorene-d10	15.228	176	342529	25.501	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	55122m	23.665	ng/ul	0.00
73) Anthracene-d10	17.098	188	532170	27.875	ng/ul	0.00
81) Pyrene-d10	19.363	212	577345	28.780	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.274	264	592786	27.965	ng/ul	0.00
Target Compounds						Qvalue

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

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