Quantitation Report (Qedit)

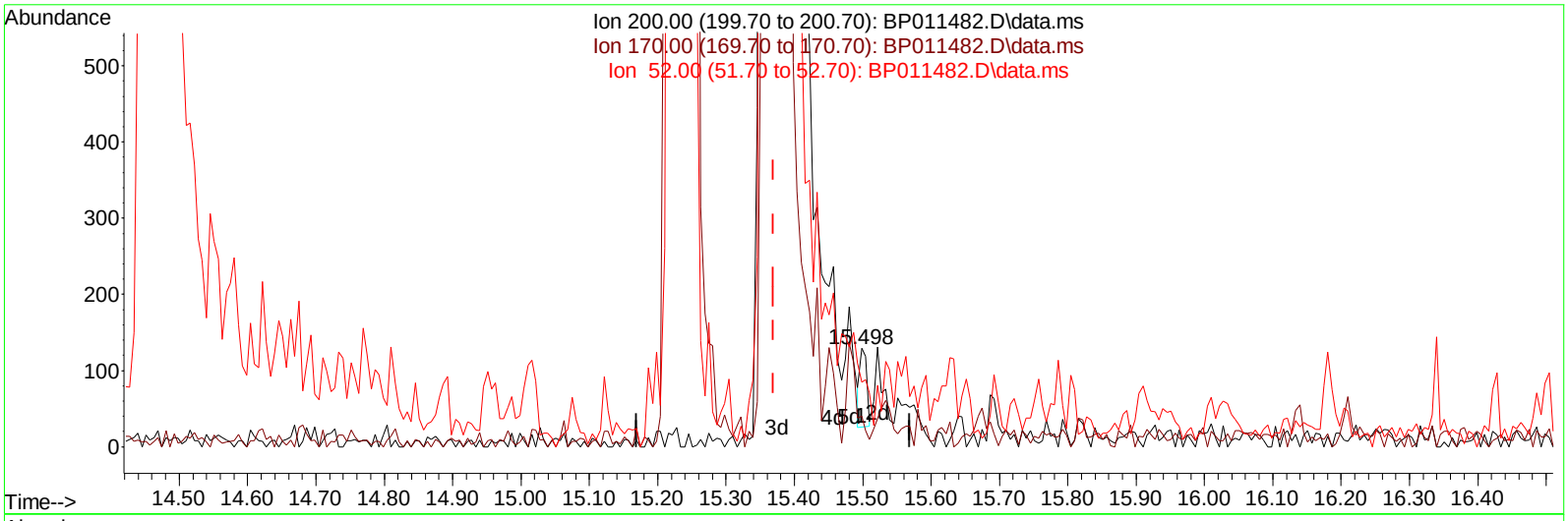
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
Data File : BP011482.D
Acq On : 18 Aug 2022 15:12
Operator : CG/JU
Sample : PB147098BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

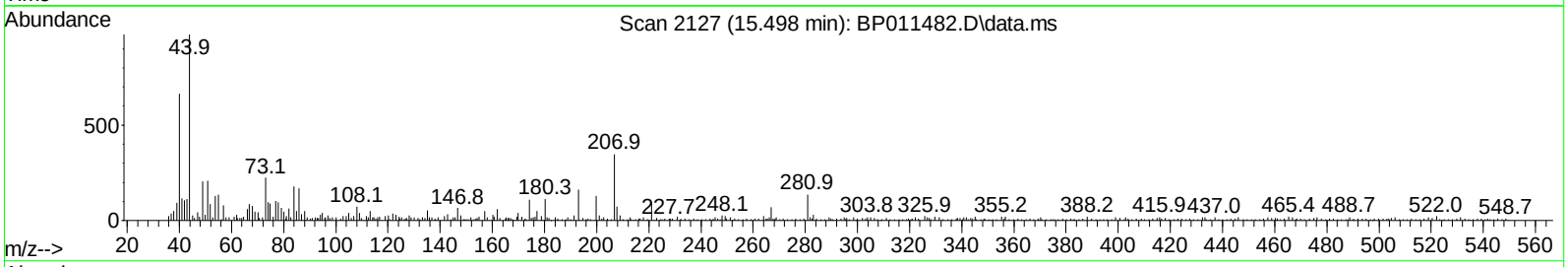
Quant Time: Aug 18 23:16:30 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 01:58:37 2022
Response via : Initial Calibration

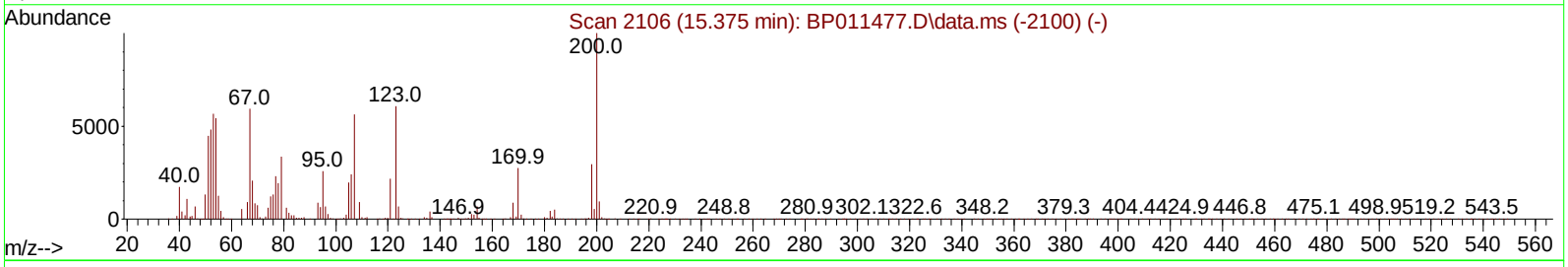
Instrument :
BNA_P
ClientSampleId :
SBLK098

Manual IntegrationsAPPROVED

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







TIC: BP011482.D\data.ms

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

15.498min (+ 0.129) 0.03 ng/ul

response 72

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	31.01
52.00	67.40	65.89
0.00	0.00	0.00

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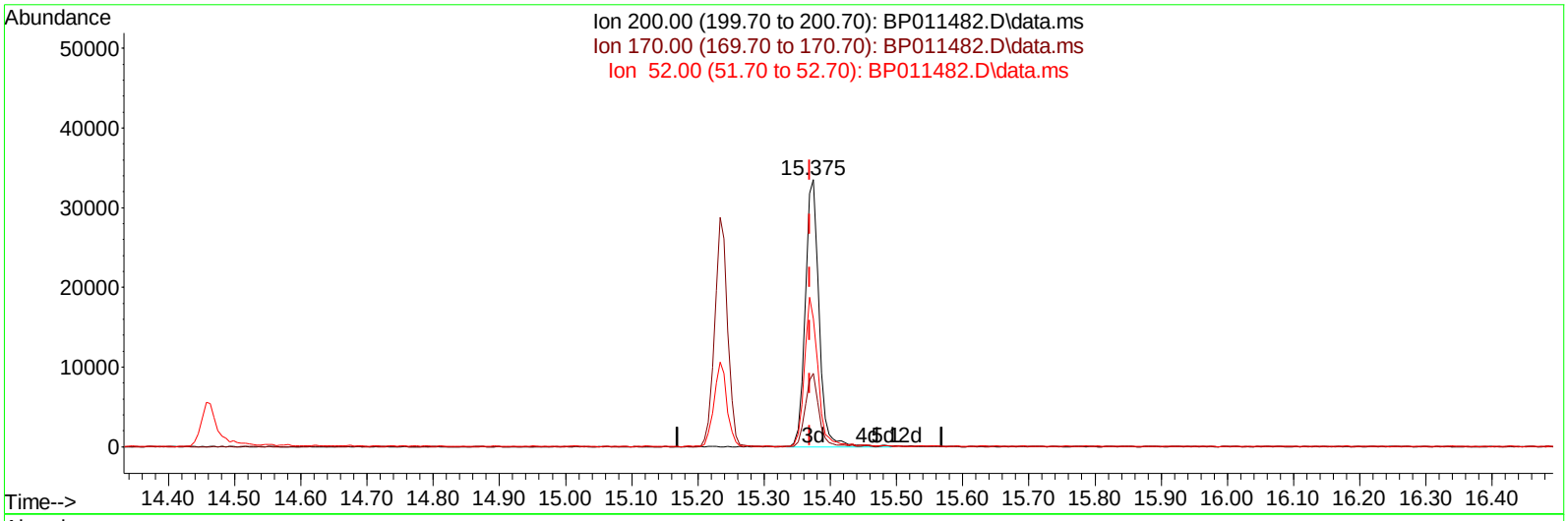

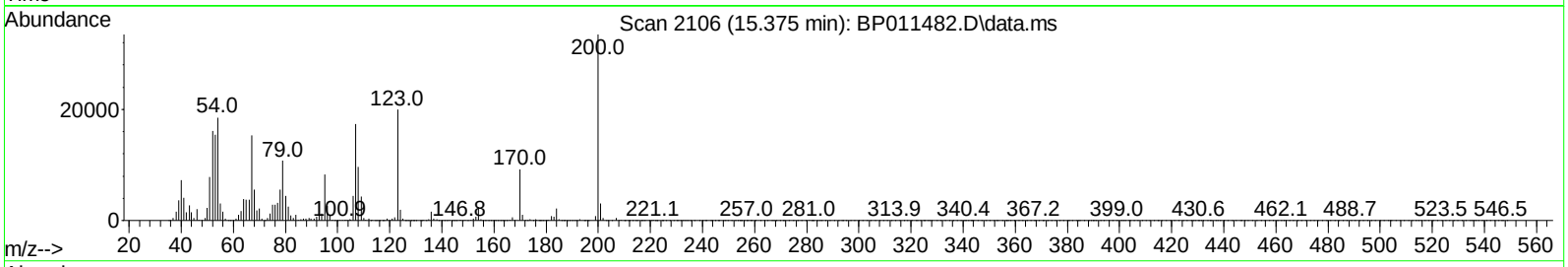

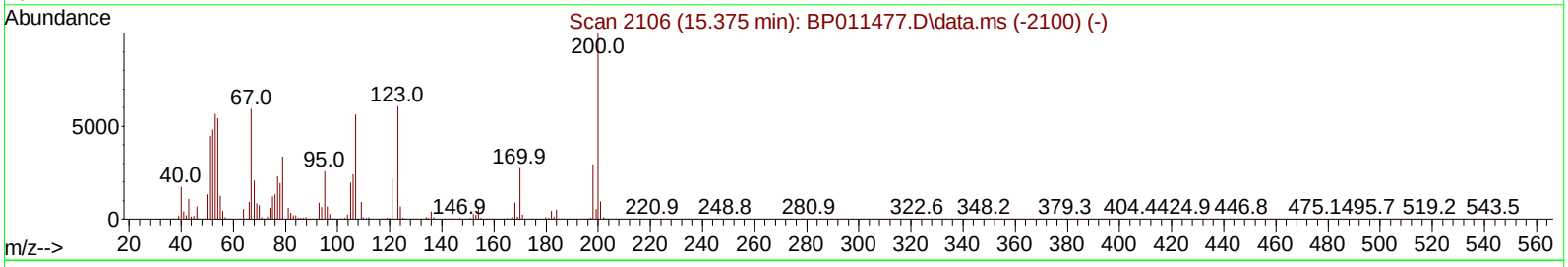

TIC: BP011482.D\data.ms

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

15.375min (+ 0.006) 23.13 ng/ul m

response 48227

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	26.30	27.53
52.00	67.40	48.10#
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	7.557	152	90279	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	368245	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	200036	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	385916	20.000	ng/ul	0.00
79) Chrysene-d12	21.133	240	366962	20.000	ng/ul	0.00
88) Perylene-d12	23.433	264	416988	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	14328	4.925	ng/uL	0.00
4) Pyridine-d5	3.475	84	174174	23.384	ng/ul	0.00
7) Phenol-d5	6.740	99	195539	23.570	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	130190	24.154	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	159139	26.080	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	151273	22.781	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	72118	27.203	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	74899	27.760	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	124682	25.035	ng/ul	0.00
31) 4-Chloroaniline-d4	10.498	131	190457	22.973	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	383655	27.117	ng/ul	0.00
49) Acenaphthylene-d8	13.922	160	406714	25.865	ng/ul	0.00
54) 4-Nitrophenol-d4	14.463	143	52171	19.139	ng/ul	0.00
60) Fluorene-d10	15.233	176	300161	25.218	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	48227m	23.127	ng/ul	0.00
73) Anthracene-d10	17.104	188	466606	27.300	ng/ul	0.00
81) Pyrene-d10	19.369	212	518189	27.910	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.286	264	559442	27.581	ng/ul	0.00
Target Compounds						Qvalue

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

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