Quantitation Report (Qedit)

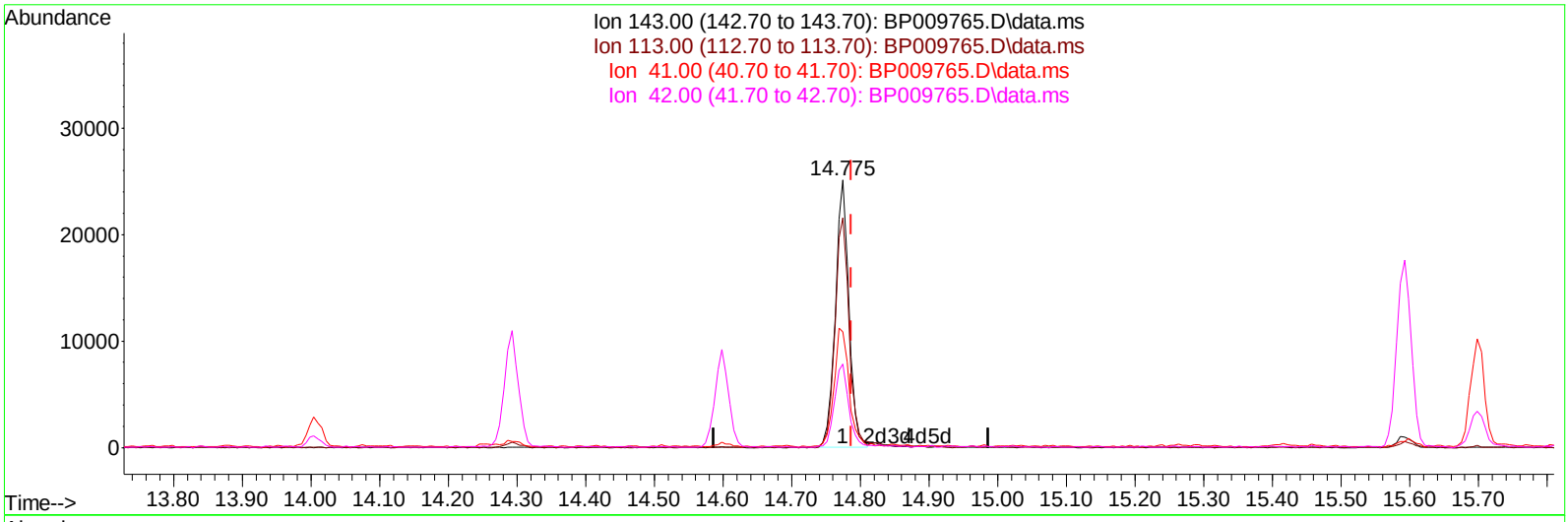
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009765.D
Acq On : 11 Apr 2022 23:16
Operator : CG/JU
Sample : N2338-15
Misc :
ALS Vial : 22 Sample Multiplier: 1

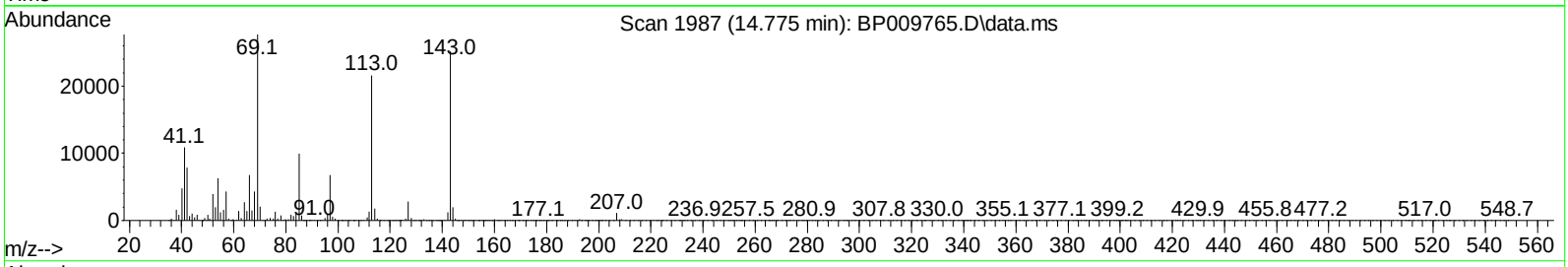
Quant Time: Apr 12 01:54:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 07 19:16:52 2022
Response via : Initial Calibration

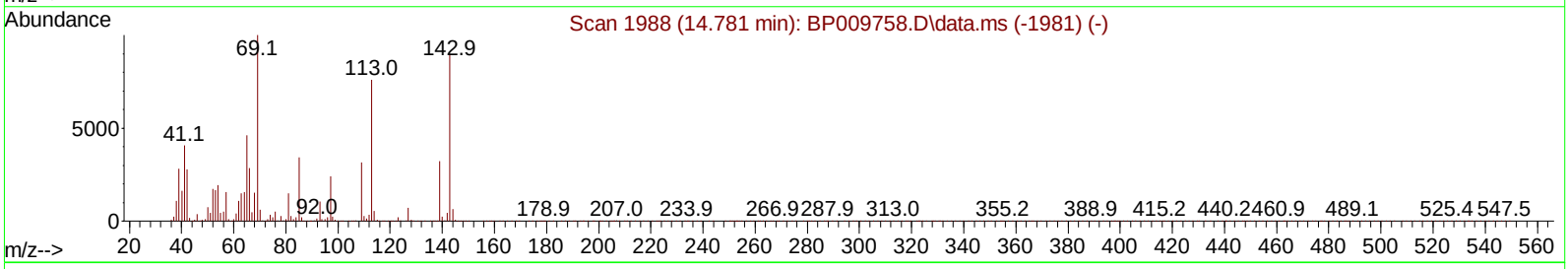
Instrument :
BNA_P
ClientSampleId :
C0J87

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/13/2022







TIC: BP009765.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 6.97 ng/ul

response 34511

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.98#
41.00	23.20	43.25#
42.00	18.00	31.28#

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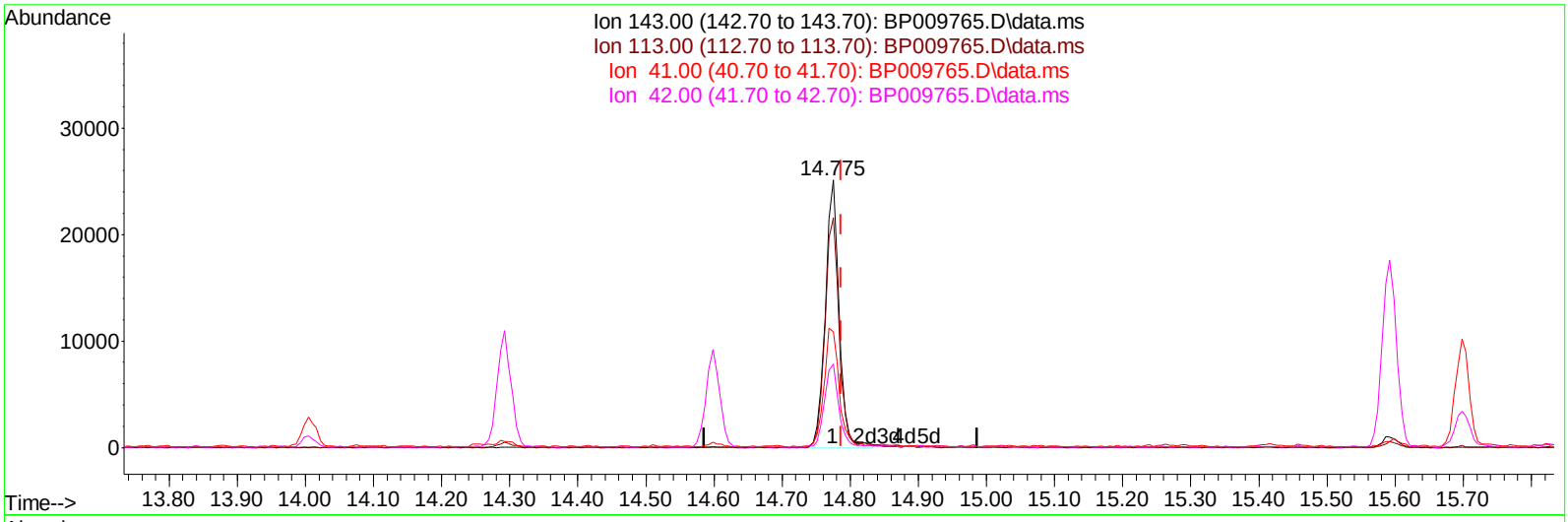

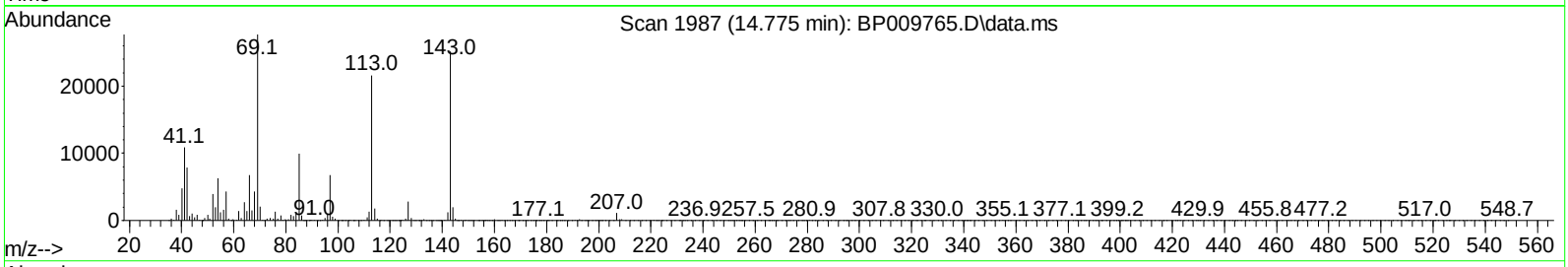

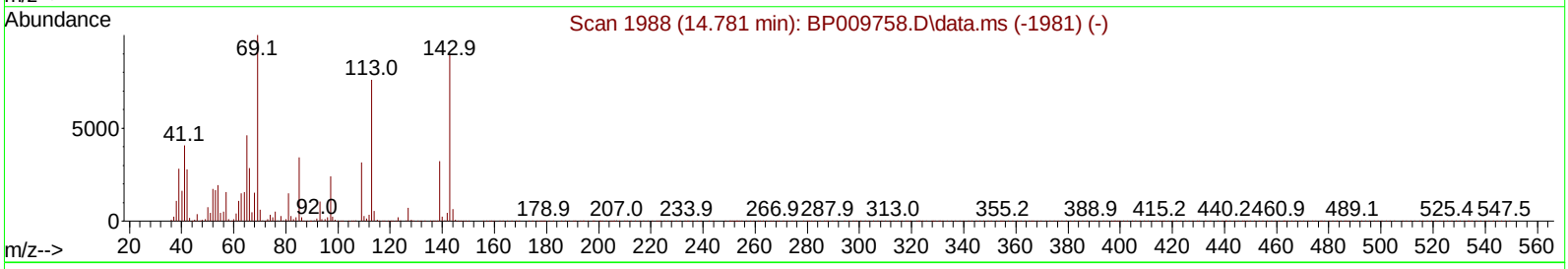

TIC: BP009765.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 7.07 ng/ul m

response 35012

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.98#
41.00	23.20	43.25#
42.00	18.00	31.28#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	107633	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	482071	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.598	164	347808	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	789231	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.427	240	864982	20.000	ng/ul	#-0.02
88) Perylene-d12	23.898	264	921201	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	13124	5.442	ng/uL	0.00
4) Pyridine-d5	3.805	84	48698	7.324	ng/ul	0.00
7) Phenol-d5	7.122	99	66253	7.011	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	193480	34.393	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	178082	24.984	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	130455	16.538	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	125415	36.077	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.851	143	130038	34.119	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	225857	28.319	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.904	131	327527	28.834	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	1010363	36.982	ng/ul	-0.01
49) Acenaphthylene-d8	14.292	160	1104556	34.126	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	35012m	7.066	ng/ul	-0.01
60) Fluorene-d10	15.592	176	875963	38.113	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	151782	32.464	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1498484	39.655	ng/ul	-0.01
81) Pyrene-d10	19.674	212	1835113	38.846	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.745	264	1964995	40.961	ng/ul	-0.02

Target Compounds Qvalue

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

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