Quantitation Report (Qedit)

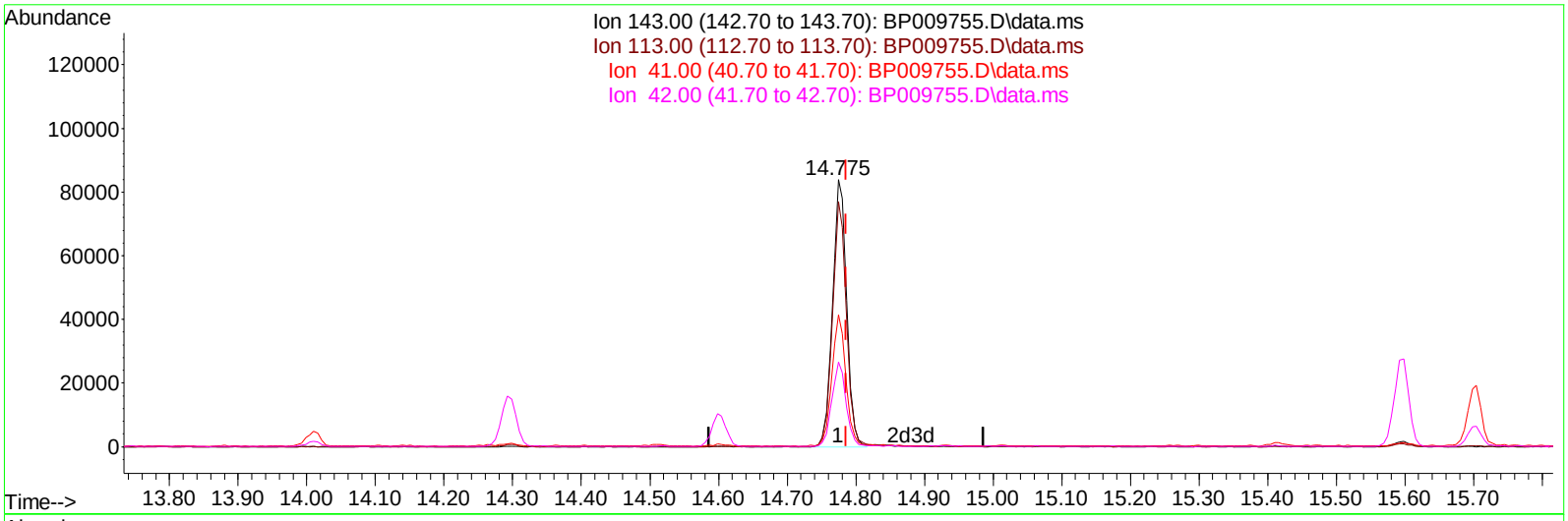
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009755.D
Acq On : 11 Apr 2022 17:04
Operator : CG/JU
Sample : N2338-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

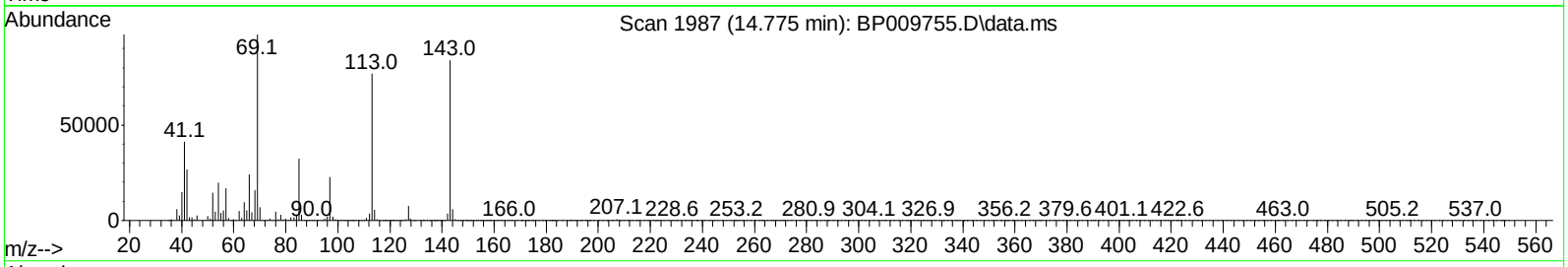
Quant Time: Apr 12 01:46:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 07 19:16:52 2022
Response via : Initial Calibration

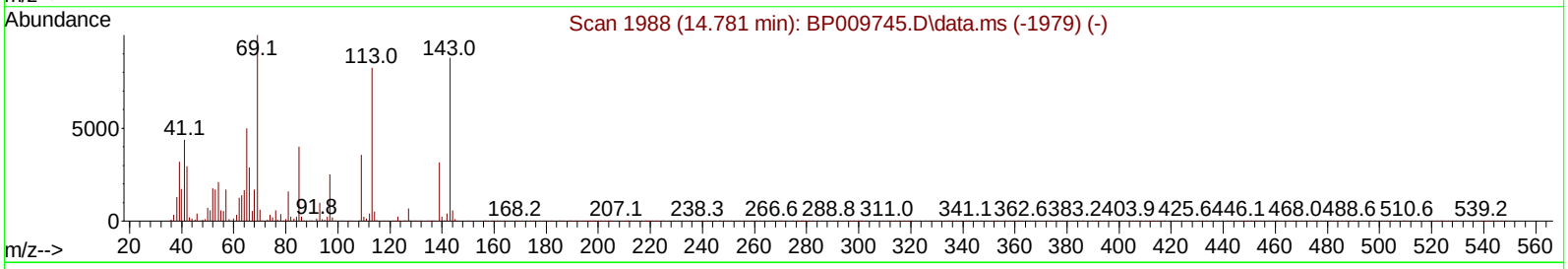
Instrument :
BNA_P
ClientSampleId :
C0J98

Manual IntegrationsAPPROVED

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







TIC: BP009755.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 20.13 ng/ul

response 119847

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	91.76#
41.00	23.20	49.29#
42.00	18.00	31.84#

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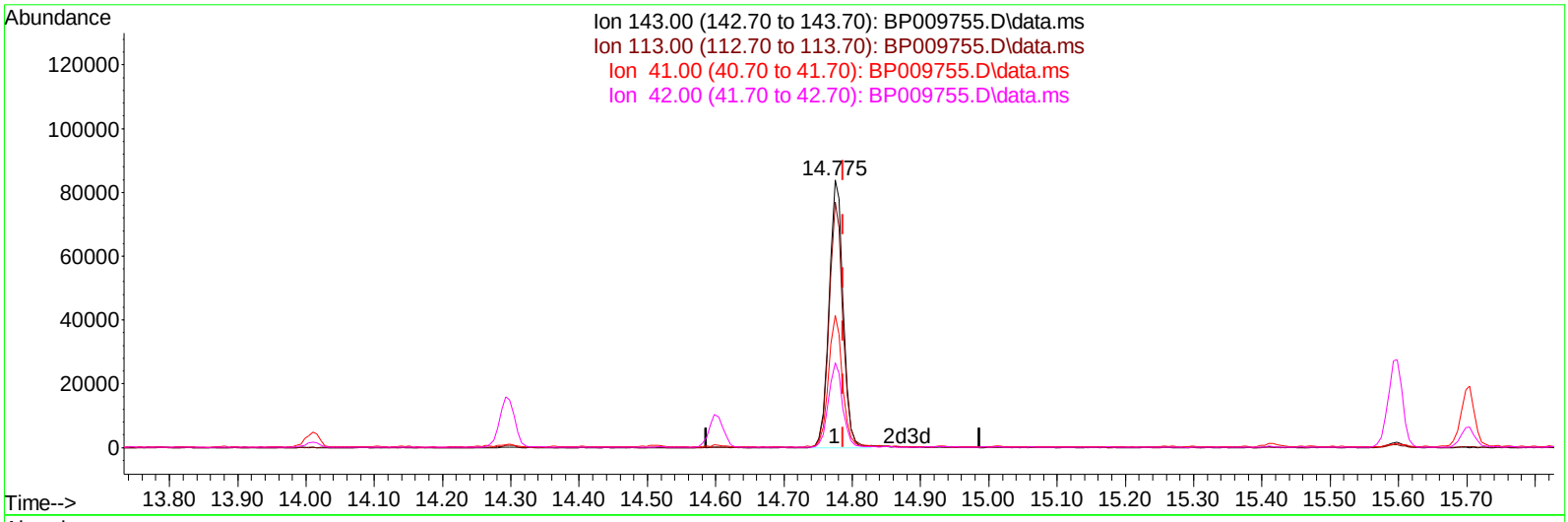

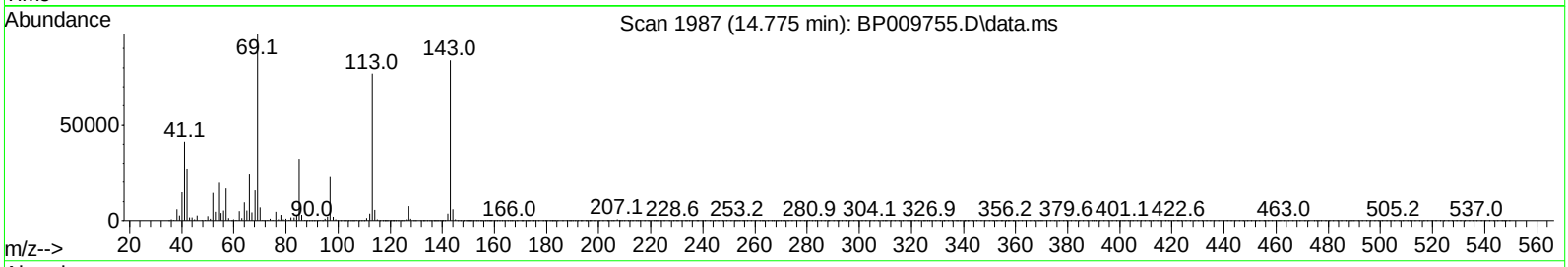

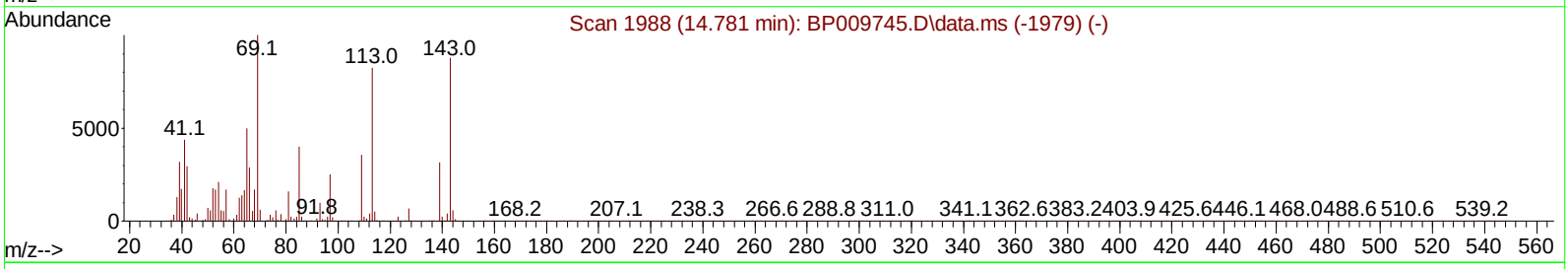

TIC: BP009755.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 20.19 ng/ul m

response 120217

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	91.76#
41.00	23.20	49.29#
42.00	18.00	31.84#

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	130076	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	576127	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	417960	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	910238	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	889720	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	772641	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	20289	6.961	ng/uL	0.00
4) Pyridine-d5	3.805	84	182221	22.676	ng/ul	0.00
7) Phenol-d5	7.122	99	200355	17.544	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	316309	46.525	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	322491	37.437	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	271919	28.523	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	209904	50.523	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	222397	48.826	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	394628	41.403	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	541377	39.879	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1700365	51.791	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	1781430	45.800	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	120217m	20.190	ng/ul	-0.01
60) Fluorene-d10	15.598	176	1409552	51.036	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	286329	53.100	ng/ul	0.00
73) Anthracene-d10	17.457	188	2348355	53.884	ng/ul	0.00
81) Pyrene-d10	19.680	212	2962694	60.970	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	2376794	59.071	ng/ul	-0.01
Target Compounds						Qvalue

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

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