Quantitation Report (Qedit)

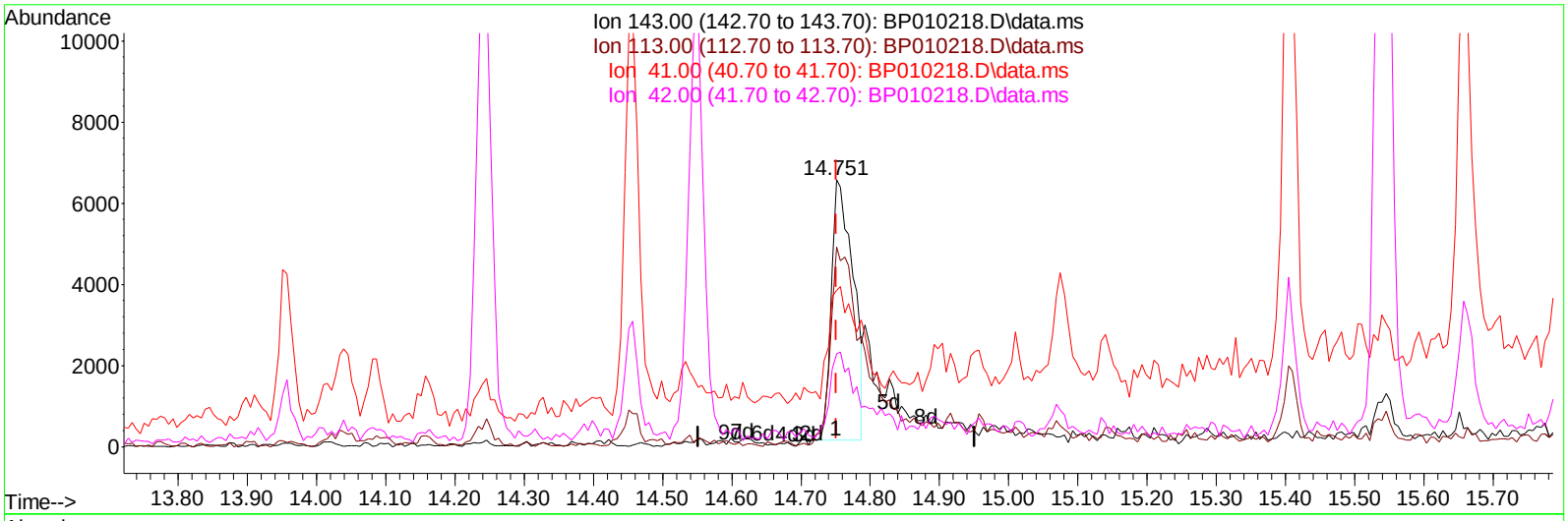
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010218.D
Acq On : 03 May 2022 21:46
Operator : CG/JU
Sample : N2670-04
Misc :
ALS Vial : 55 Sample Multiplier: 1

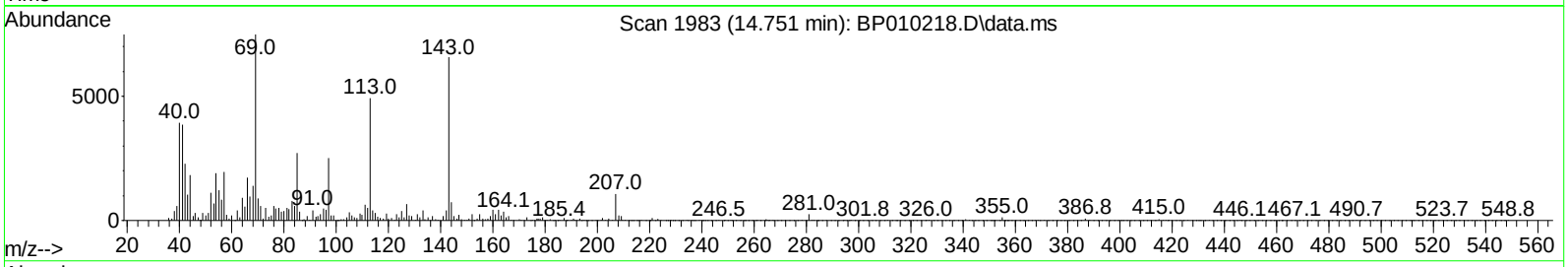
Quant Time: May 04 01:25:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

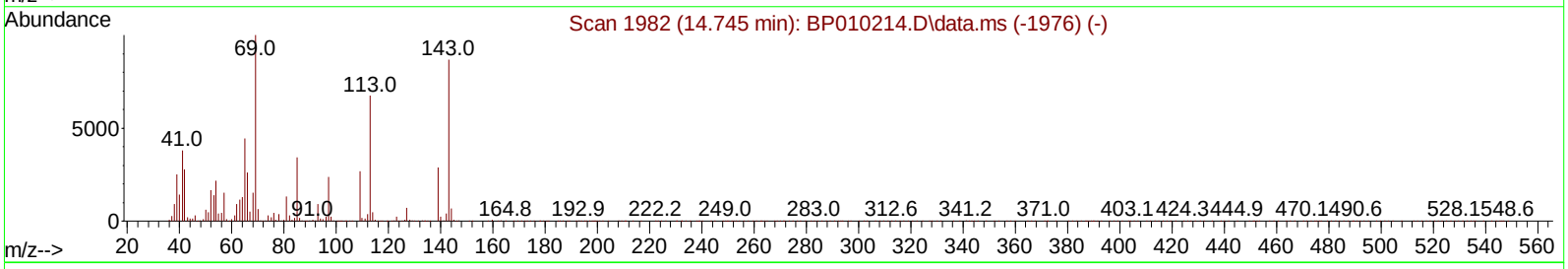
Instrument :
BNA_P
ClientSampleId :
C0W06

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/04/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010218.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 2.12 ng/ul

response 13926

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.87#
41.00	23.20	58.49#
42.00	18.00	34.89#

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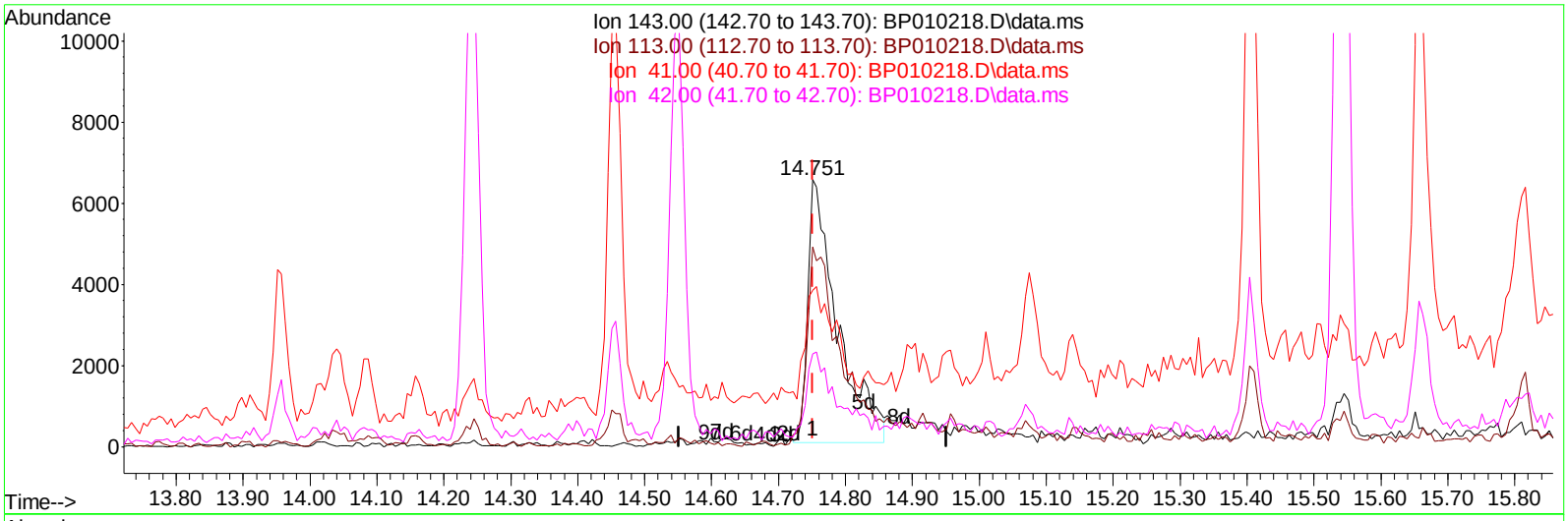

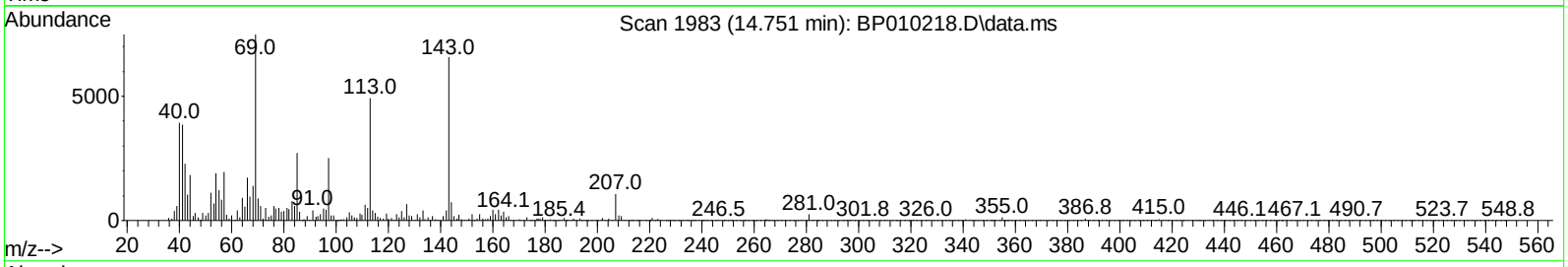

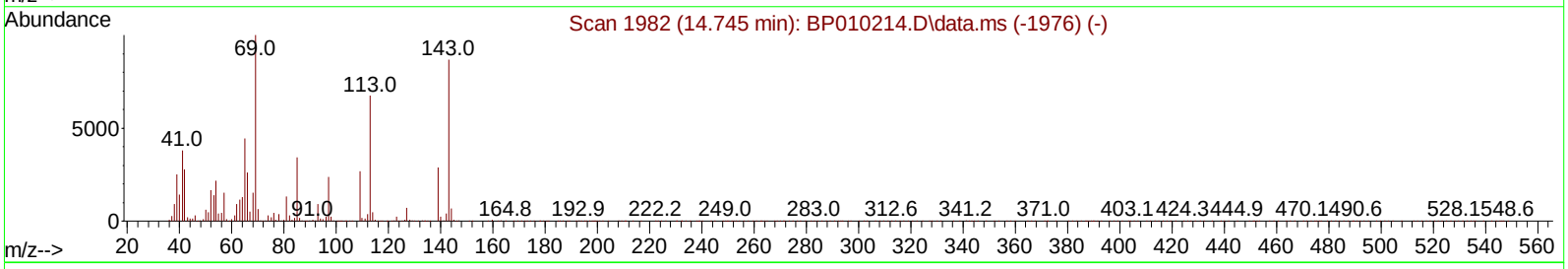

TIC: BP010218.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 3.03 ng/ul m

response 19957

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.87#
41.00	23.20	58.49#
42.00	18.00	34.89#

Quantitation Report (QT Reviewed)

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Supervised By :Yogesh Patel 05/05/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	166697	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	717139	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	459276	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	865451	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	712325	20.000	ng/ul	#-0.01
88) Perylene-d12	23.827	264	685485	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	19085	4.992	ng/uL	0.00
4) Pyridine-d5	3.781	84	82809	7.700	ng/ul	0.00
7) Phenol-d5	7.081	99	79950	5.273	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	304194	33.434	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	261144	23.256	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	166565	13.259	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	192127	33.684	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	192694	30.934	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	308073	25.404	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	394994	22.717	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1245970	34.875	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1411160	32.636	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143	19957m	3.034	ng/ul	0.00
60) Fluorene-d10	15.545	176	1051479	34.214	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	141979	25.945	ng/ul	0.00
73) Anthracene-d10	17.404	188	1552217	37.073	ng/ul	0.00
81) Pyrene-d10	19.633	212	1664116	41.079	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1428284	39.744	ng/ul	-0.01
Target Compounds						Qvalue

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

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Supervised By :Yogesh Patel 05/05/2022