Quantitation Report (Qedit)

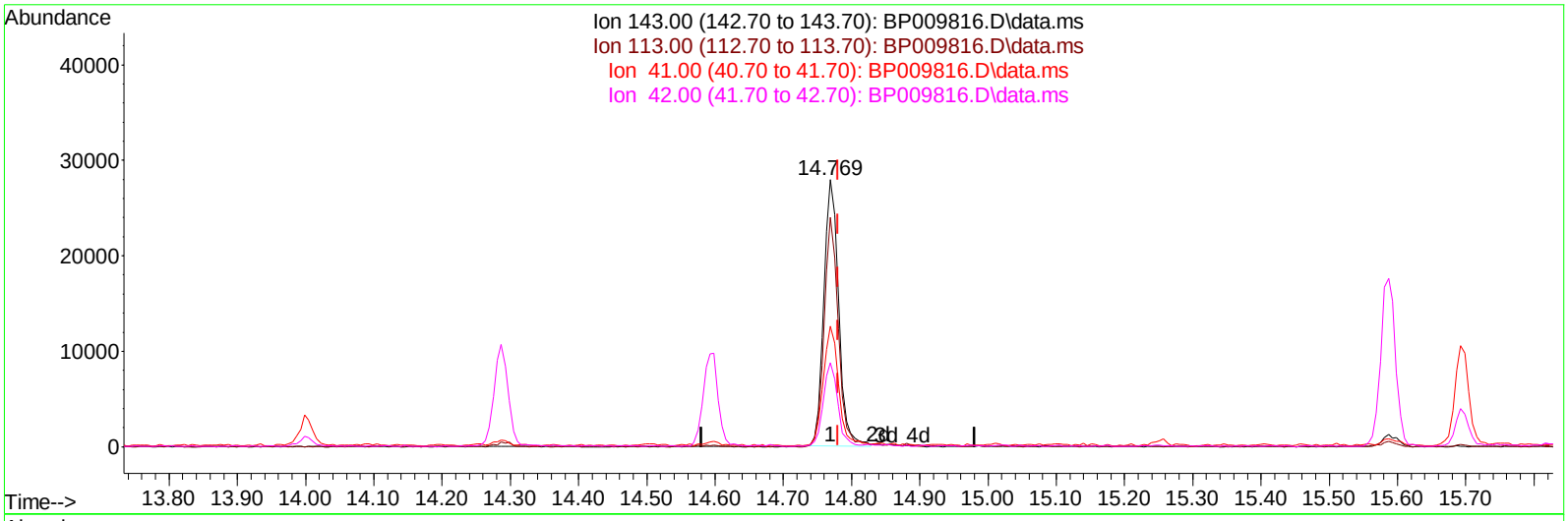
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041322\
Data File : BP009816.D
Acq On : 13 Apr 2022 14:06
Operator : CG/JU
Sample : N2342-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

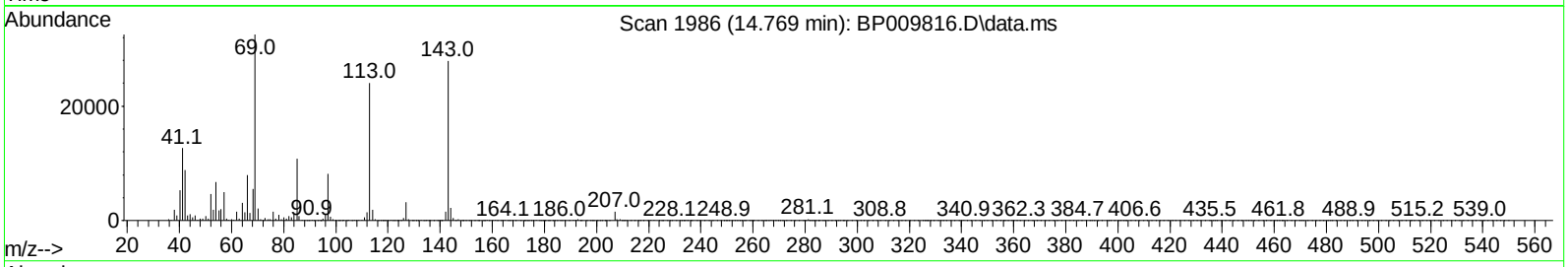
Quant Time: Apr 14 01:06:09 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

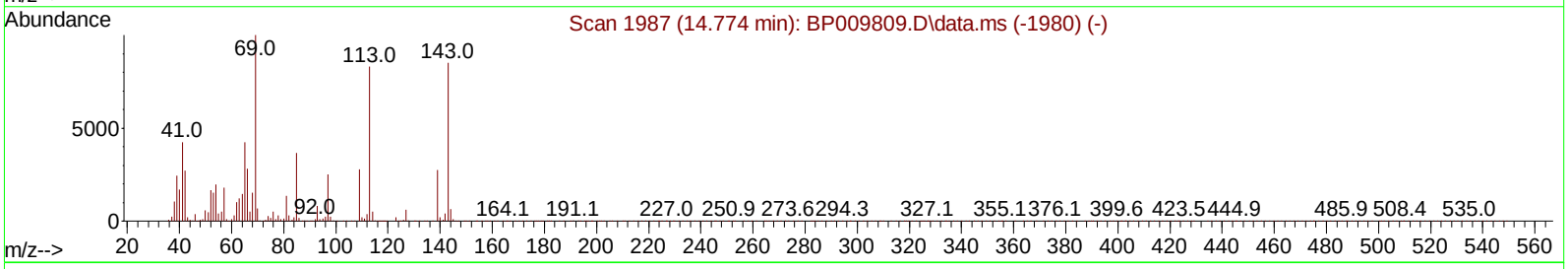
Instrument :
BNA_P
ClientSampleId :
C0HY3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/14/2022
Supervised By :Yogesh Patel 04/14/2022







TIC: BP009816.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 6.85 ng/ul

response 41858

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.04#
41.00	23.20	45.27#
42.00	18.00	31.55#

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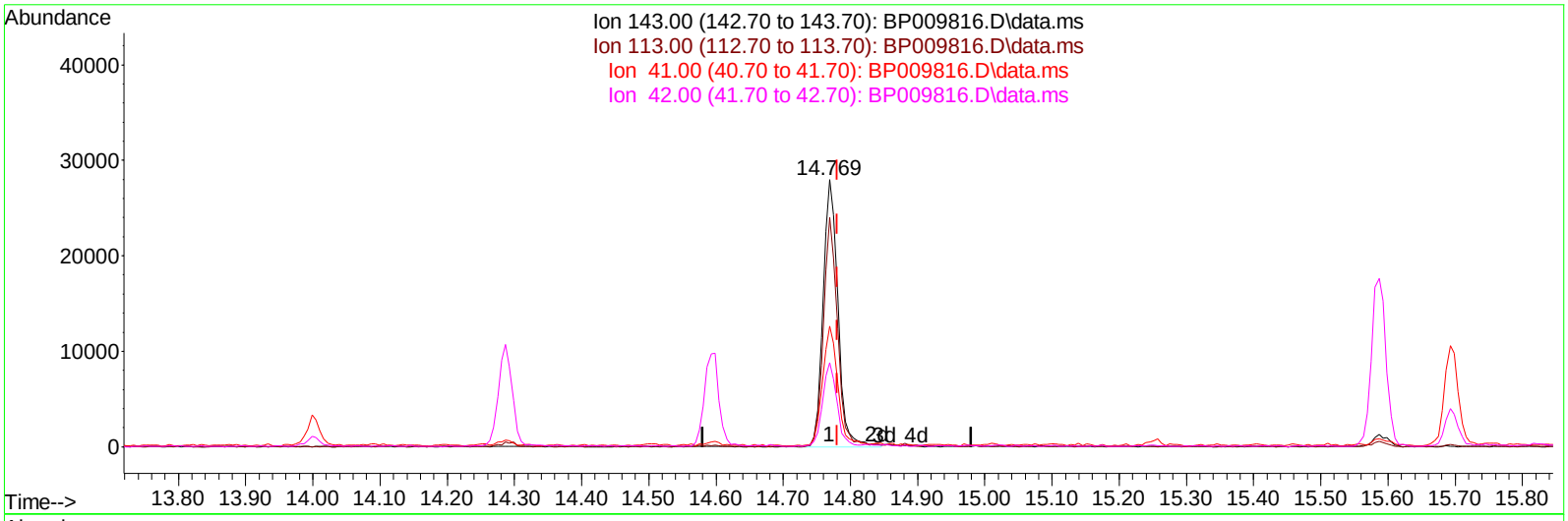

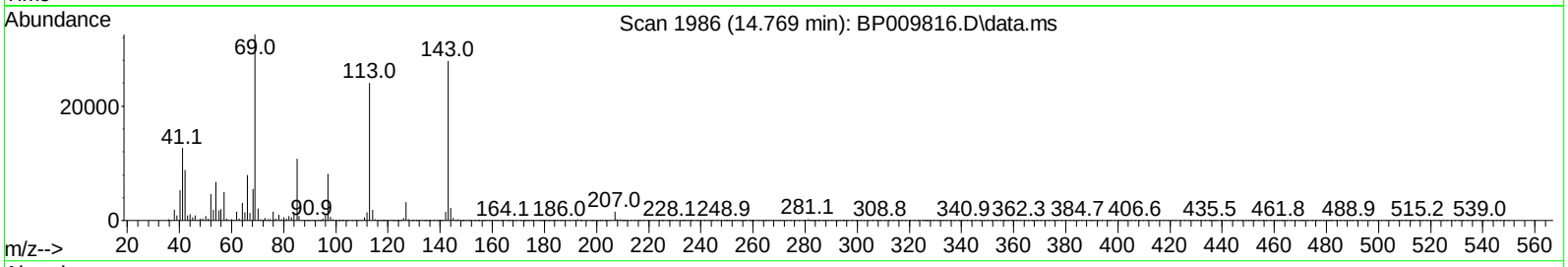

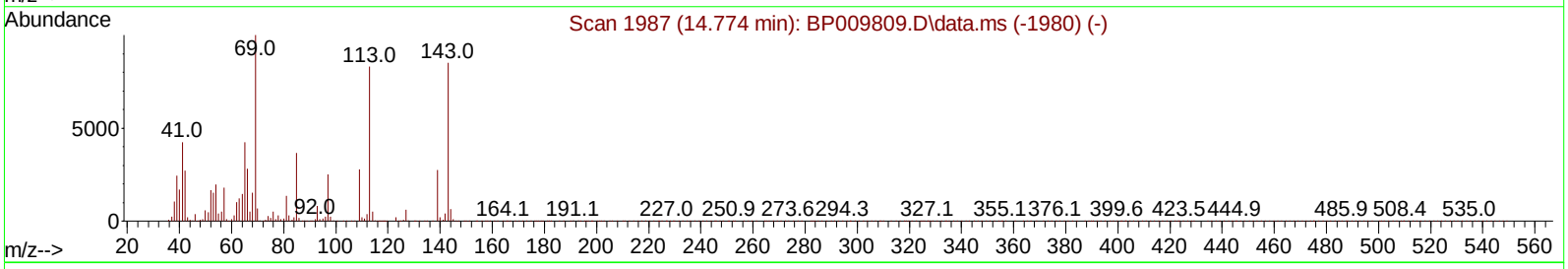

TIC: BP009816.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 6.96 ng/ul m

response 42571

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.04#
41.00	23.20	45.27#
42.00	18.00	31.55#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.964	152	132824	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	600588	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	429254	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	952985	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.427	240	958366	20.000	ng/ul	# 0.00
88) Perylene-d12	23.892	264	830051	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	16693	5.609	ng/uL	0.00
4) Pyridine-d5	3.805	84	82050	9.999	ng/ul	0.00
7) Phenol-d5	7.117	99	86483	7.416	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	222121	31.996	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	215602	24.511	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	164057	16.853	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	144737	33.419	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	149507	31.486	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	258628	26.029	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	368548	26.042	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1082590	32.107	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1181834	29.585	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	42571m	6.962	ng/ul	-0.01
60) Fluorene-d10	15.587	176	930731	32.812	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	163908	29.033	ng/ul	0.00
73) Anthracene-d10	17.445	188	1602059	35.111	ng/ul	0.00
81) Pyrene-d10	19.675	212	1960921	37.464	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1592557	36.842	ng/ul	0.00

Target Compounds Qvalue

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

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