Quantitation Report (Qedit)

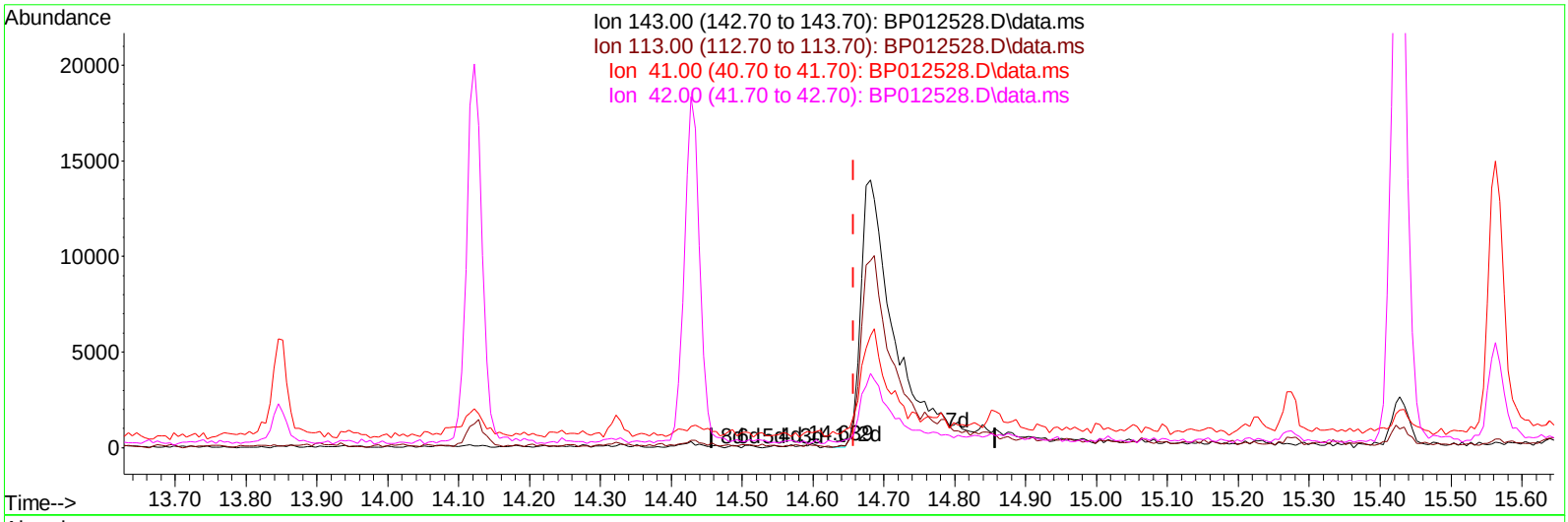
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
Data File : BP012528.D
Acq On : 07 Nov 2022 13:15
Operator : CG/JU
Sample : N5353-10
Misc :
ALS Vial : 5 Sample Multiplier: 1

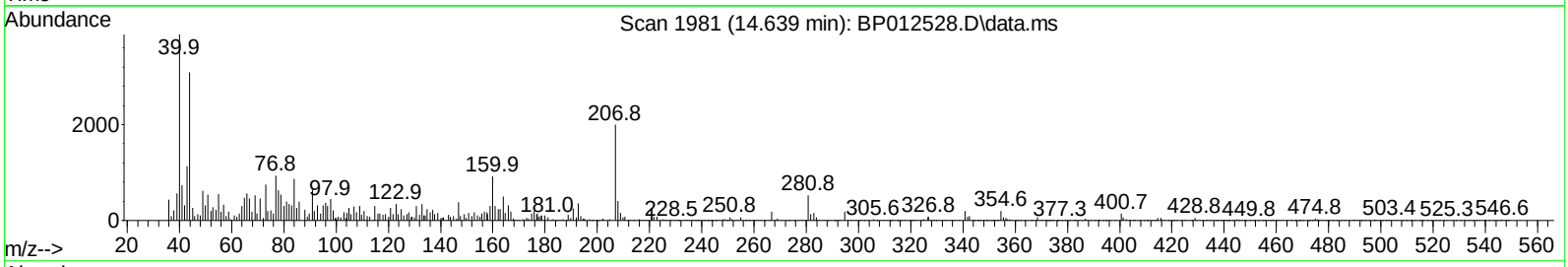
Quant Time: Nov 07 21:39:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 04 02:53:36 2022
Response via : Initial Calibration

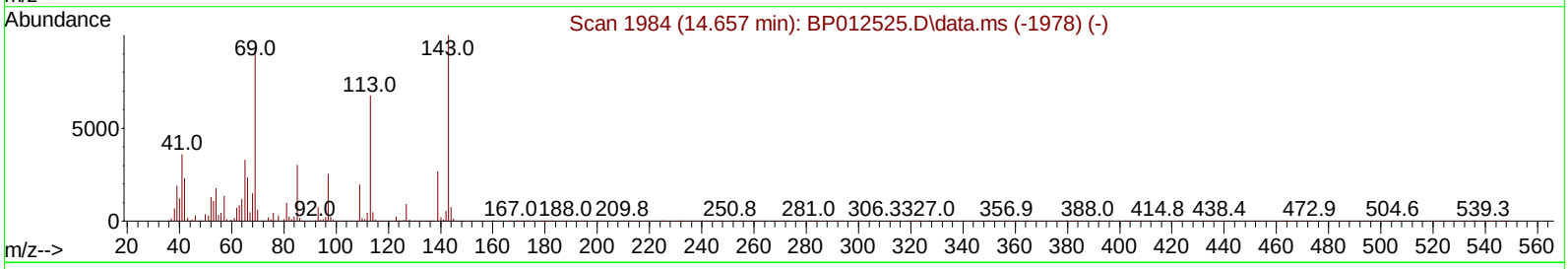
Instrument :
BNA_P
ClientSampleId :
DBWU7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/08/2022
Supervised By :mohammad ahmed 11/14/2022







TIC: BP012528.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.639min (-0.018) 0.00 ng/ul

response 58

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	61.67
41.00	34.50	617.50#
42.00	23.60	261.67#

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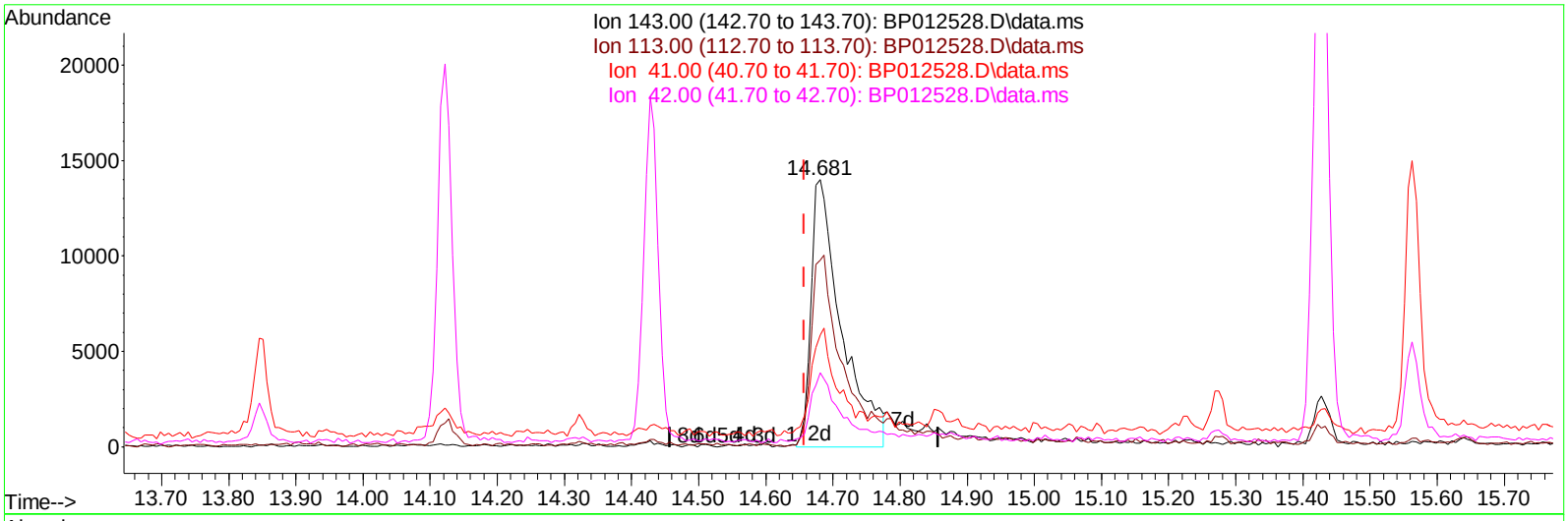

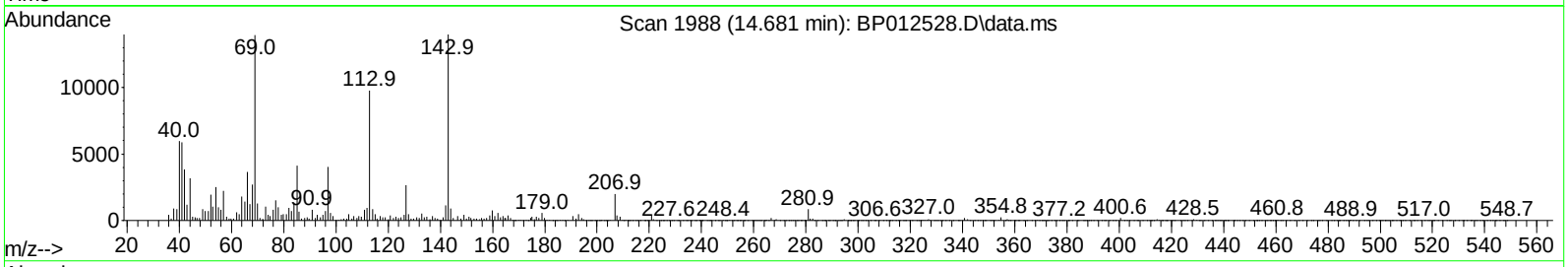

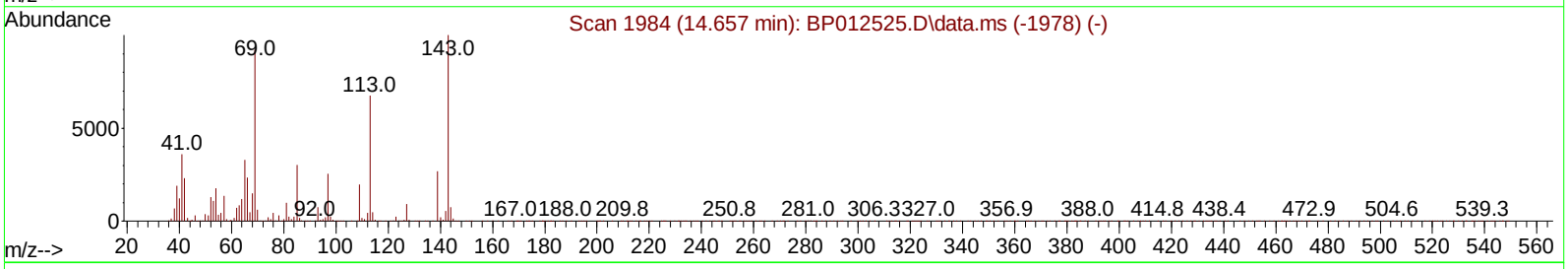

TIC: BP012528.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.023) 2.83 ng/ul m

response 44029

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	69.81
41.00	34.50	41.88#
42.00	23.60	27.61

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	371710	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1764009	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	1201343	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2704559	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	2460006	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	2283910	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	33438	3.622	ng/uL	0.00
4) Pyridine-d5	3.658	84	231426	8.727	ng/ul	0.00
7) Phenol-d5	6.957	99	200142	6.003	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	564416	28.359	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	600720	24.674	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	421514	15.986	ng/ul	0.00
21) Nitrobenzene-d5	8.946	128	410867	35.917	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	441384	37.175	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	841703	32.711	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	707643	18.470	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	2957049	35.957	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	3253689	34.621	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	44029m	2.829	ng/ul	0.02
60) Fluorene-d10	15.428	176	2610736	37.079	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	323039	22.941	ng/ul	0.00
73) Anthracene-d10	17.292	188	4206275	36.095	ng/ul	0.00
81) Pyrene-d10	19.533	212	4831336	39.117	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	4100390	36.743	ng/ul	0.00

Target Compounds Qvalue

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

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