Quantitation Report (Qedit)

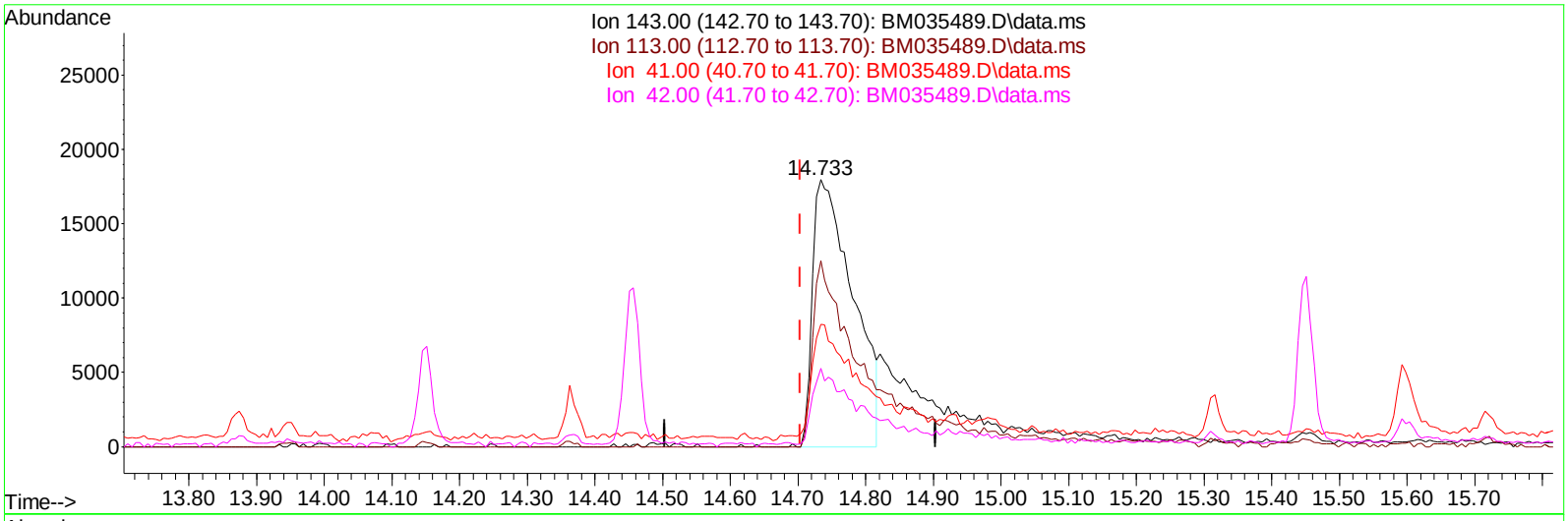
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
Data File : BM035489.D
Acq On : 17 Jun 2022 11:31
Operator : CG/JU
Sample : N3226-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

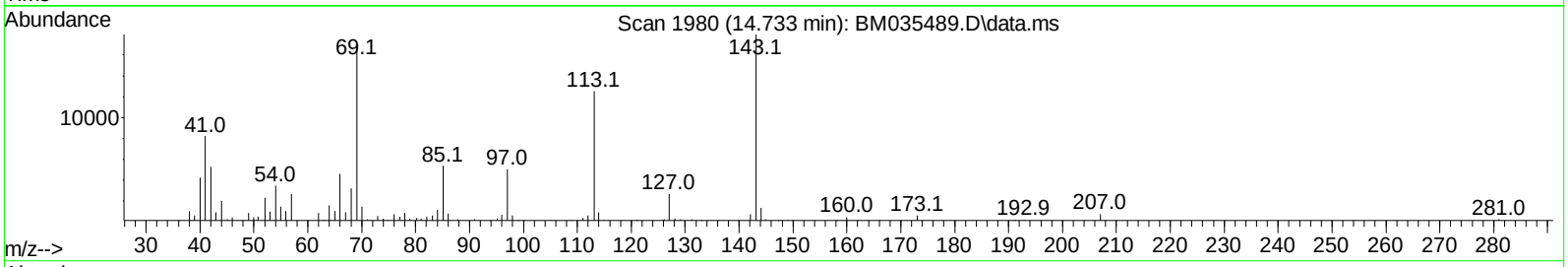
Quant Time: Jun 17 23:00:56 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 01:58:49 2022
Response via : Initial Calibration

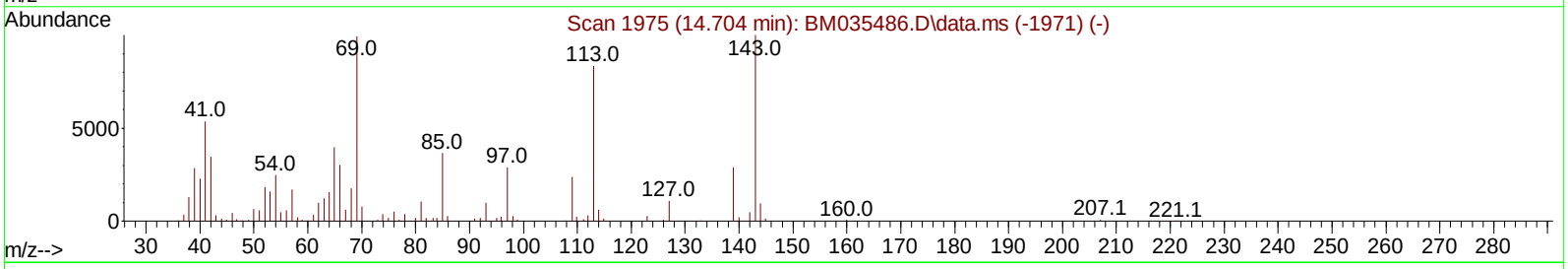
Instrument :
BNA_M
ClientSampleId :
EW4P1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/18/2022
Supervised By :Yogesh Patel 06/22/2022







TIC: BM035489.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.030) 11.89 ng/ul

response 74542

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	69.66
41.00	41.00	45.83
42.00	29.30	29.41

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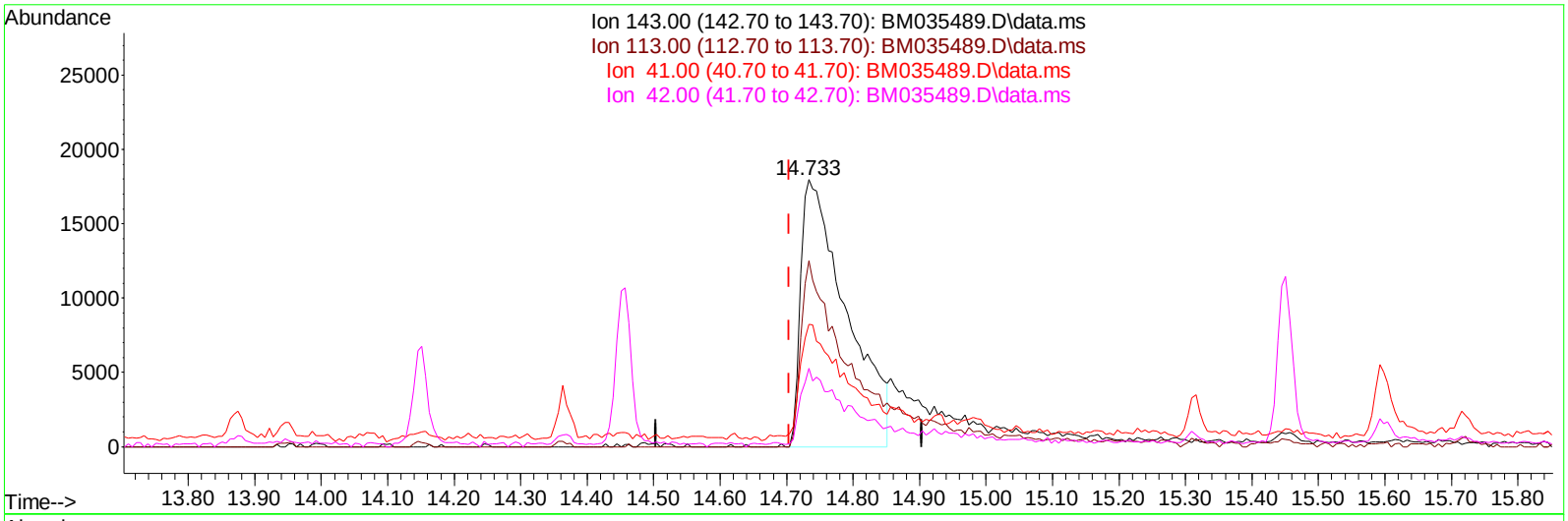

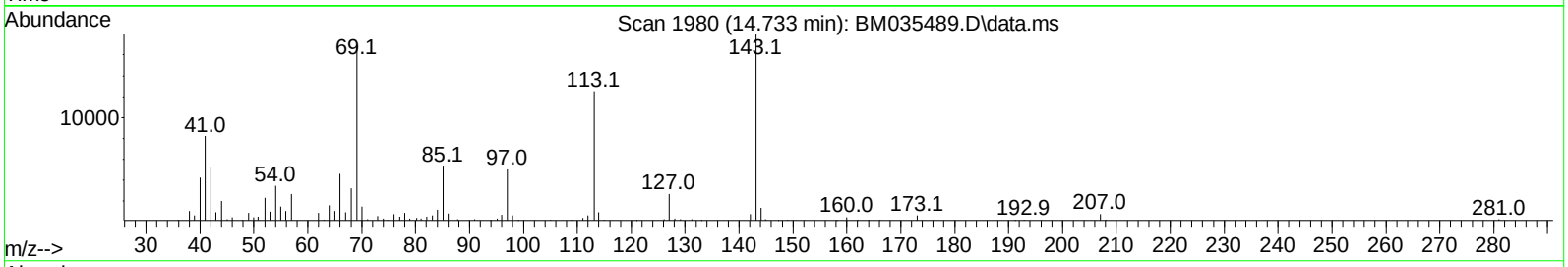

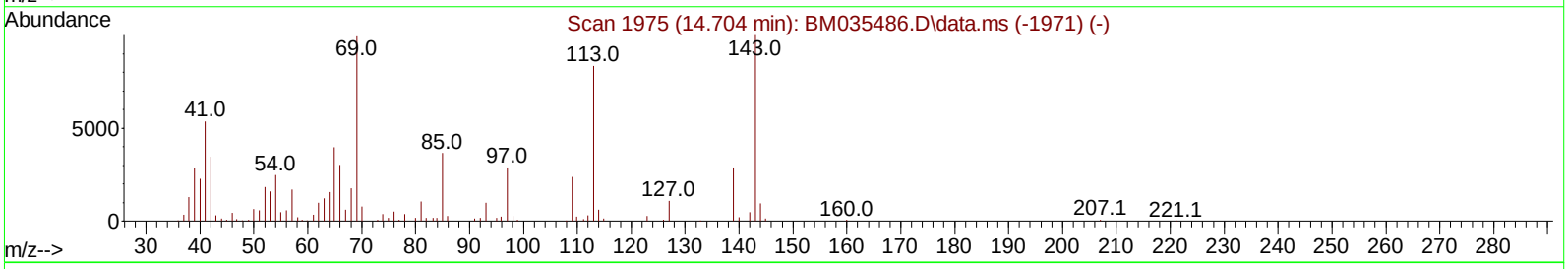

TIC: BM035489.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.030) 13.63 ng/ul m

response 85442

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	69.66
41.00	41.00	45.83
42.00	29.30	29.41

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	228184	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	979077	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	600076	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1235889	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240	1278896	20.000	ng/ul	0.00
88) Perylene-d12	23.733	264	1382272	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	18596	3.253	ng/uL	0.00
4) Pyridine-d5	3.699	84	124276	8.827	ng/ul	0.01
7) Phenol-d5	7.004	99	269904	15.434	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	178790	16.440	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	252567	16.674	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113	237195	16.202	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	115214	15.809	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	122840	15.631	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	205964	15.436	ng/ul	0.01
31) 4-Chloroaniline-d4	10.763	131	312065	14.864	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	801886	19.645	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	910312	17.997	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	85442m	13.627	ng/ul	0.03
60) Fluorene-d10	15.451	176	714328	20.211	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	93327	12.666	ng/ul	0.00
73) Anthracene-d10	17.304	188	1244311	22.399	ng/ul	0.00
81) Pyrene-d10	19.598	212	1571018	24.986	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264	1652035	24.779	ng/ul	0.00
Target Compounds						Qvalue
16) Acetophenone	8.663	105	26025	1.172	ng/ul	94

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

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