Quantitation Report (Qedit)

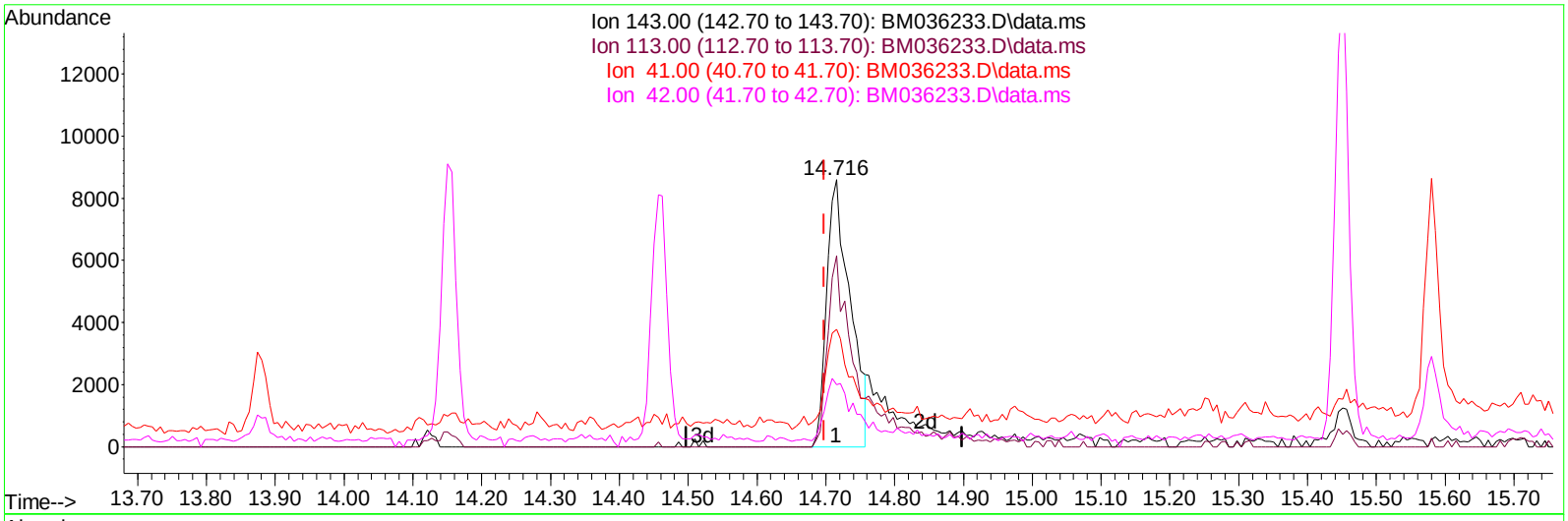
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036233.D
Acq On : 12 Aug 2022 16:55
Operator : CG/JU
Sample : N4095-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

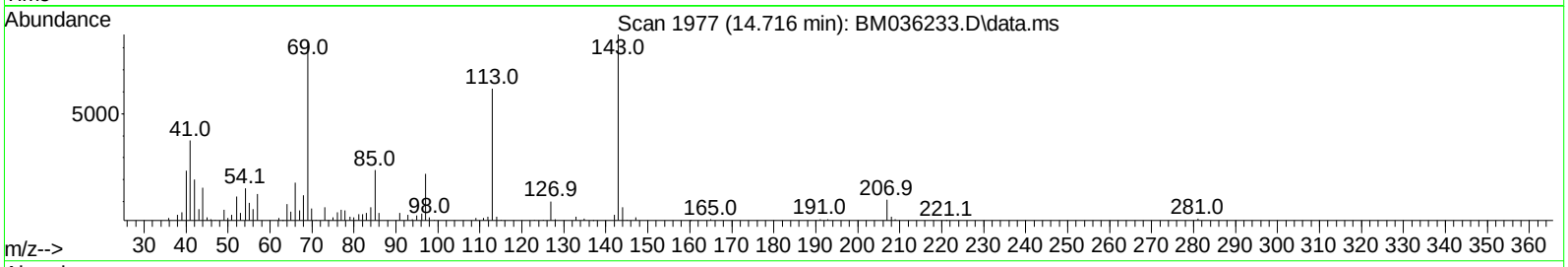
Quant Time: Aug 12 17:23:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 12 11:23:11 2022
Response via : Initial Calibration

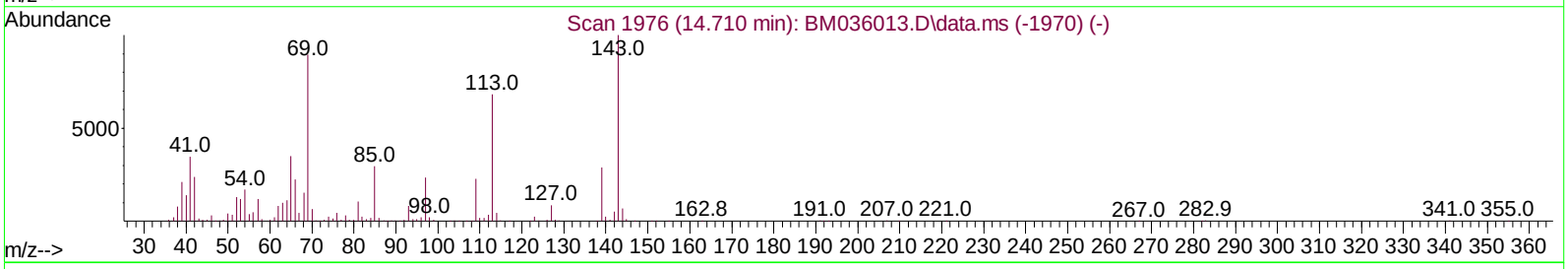
Instrument :
BNA_M
ClientSampleId :
GBJM7

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/13/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BM036233.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.018) 3.01 ng/ul

response 20243

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	71.49
41.00	41.00	44.06
42.00	29.30	23.28#

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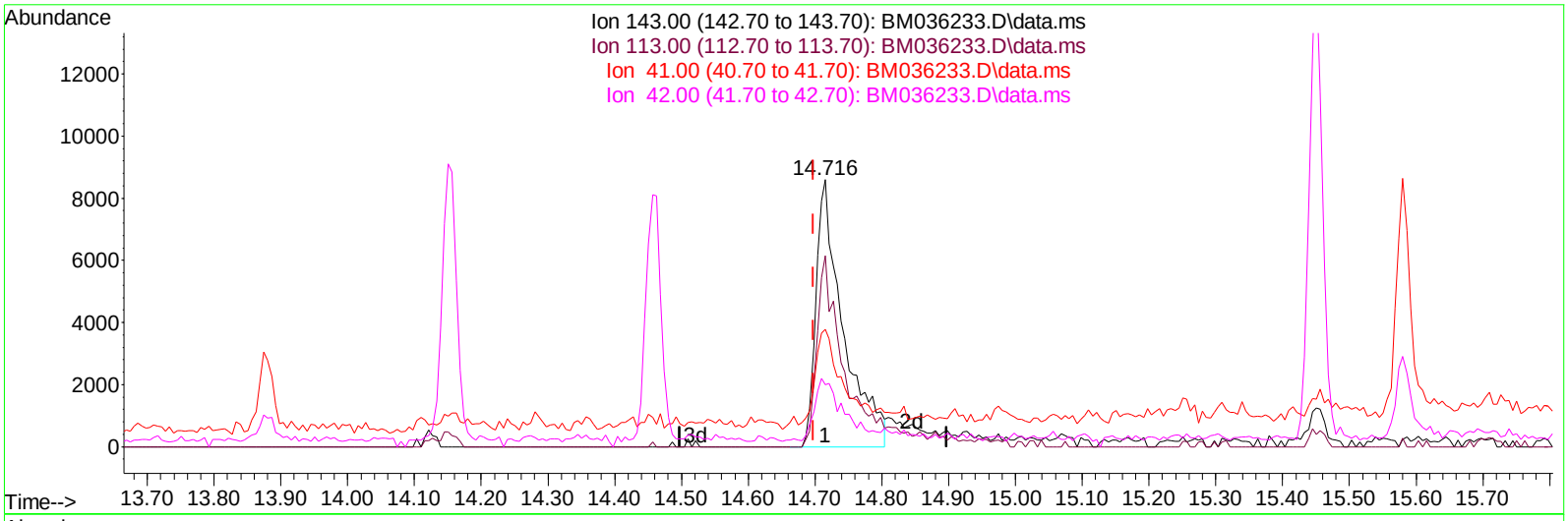

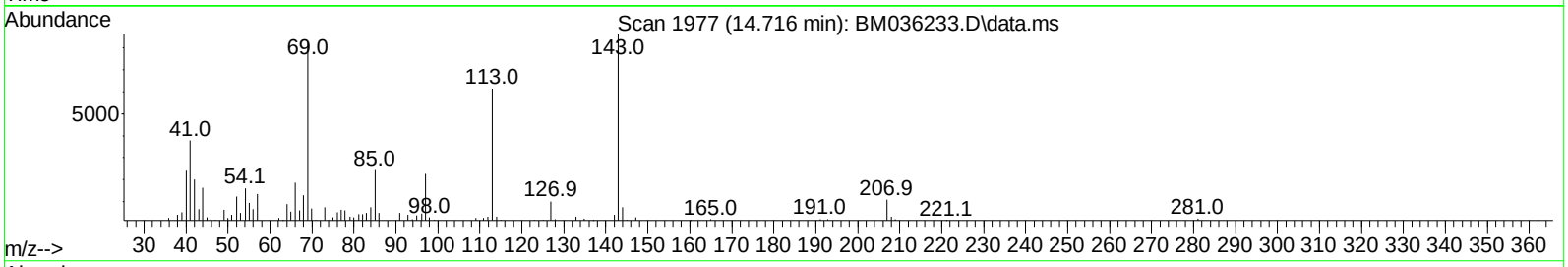

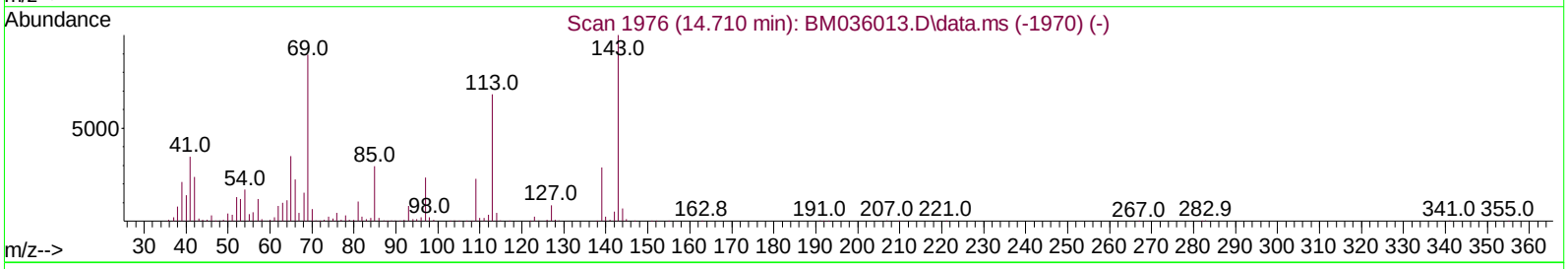

TIC: BM036233.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.018) 3.63 ng/ul m

response 24436

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	71.49
41.00	41.00	44.06
42.00	29.30	23.28#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	201934	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	837131	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	480594	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	965414	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	939734	20.000	ng/ul	0.00
88) Perylene-d12	23.727	264	1009392	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	24955	4.650	ng/uL	0.00
4) Pyridine-d5	3.663	84	39295	2.673	ng/ul	0.02
7) Phenol-d5	7.010	99	94906	5.604	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	294638	26.100	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	299097	21.279	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	174870	12.747	ng/ul	0.00
21) Nitrobenzene-d5	8.986	128	195224	29.430	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	189055	29.253	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	305203	26.111	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	454136	22.950	ng/ul	0.00
46) Dimethylphthalate-d6	13.880	166	1055419	32.239	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1332773	31.908	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	24436m	3.635	ng/ul	0.02
60) Fluorene-d10	15.451	176	979640	33.115	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	131042	25.252	ng/ul	0.00
73) Anthracene-d10	17.304	188	1599572	36.648	ng/ul	0.00
81) Pyrene-d10	19.598	212	1830849	35.153	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264	1899543	37.067	ng/ul	0.00

Target Compounds	Qvalue

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

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