Quantitation Report (Qedit)

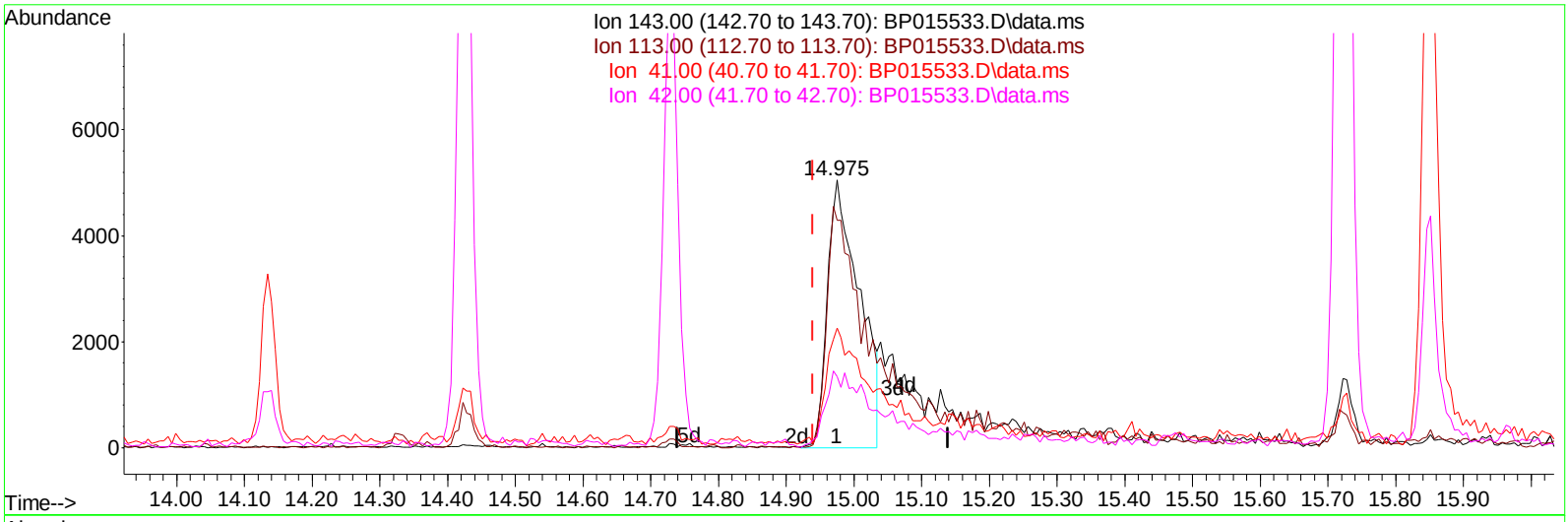
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015533.D
Acq On : 14 Jun 2023 23:56
Operator : MA/JU
Sample : 03104-05
Misc :
ALS Vial : 27 Sample Multiplier: 1

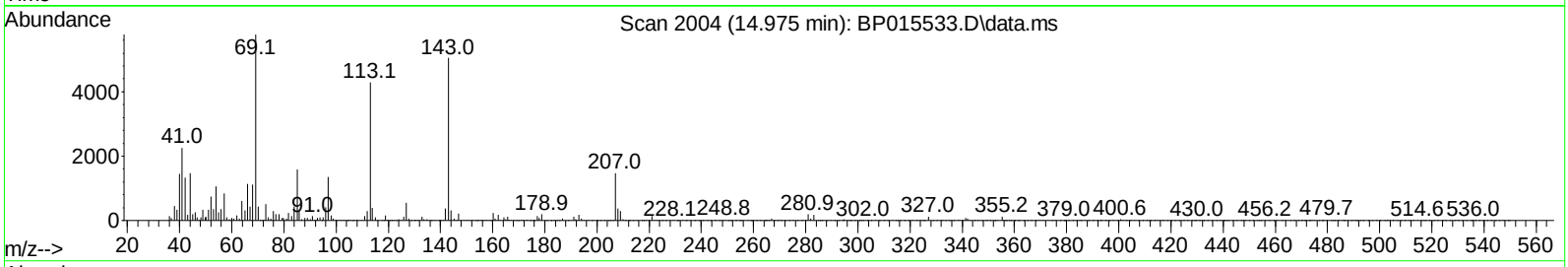
Quant Time: Jun 15 00:45:55 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

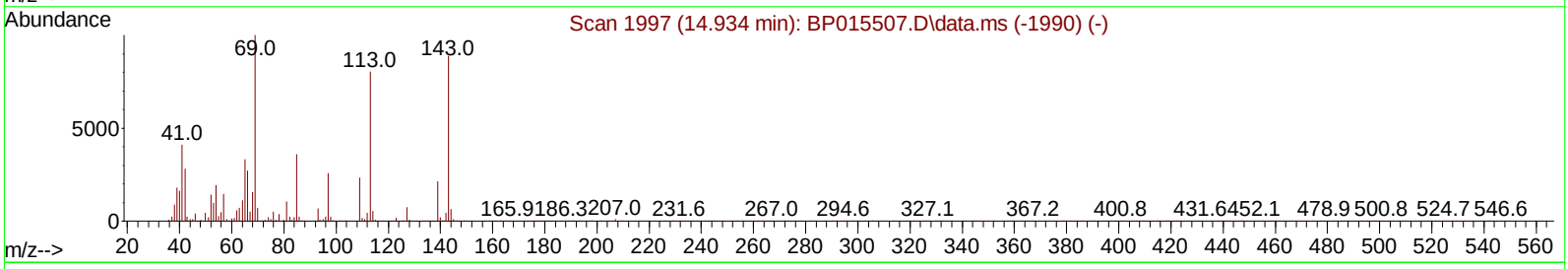
Instrument :
BNA_P
ClientSampleId :
C0NM1

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023







TIC: BP015533.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.035) 3.38 ng/ul

response 16651

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	85.01
41.00	42.20	44.70
42.00	29.40	26.48

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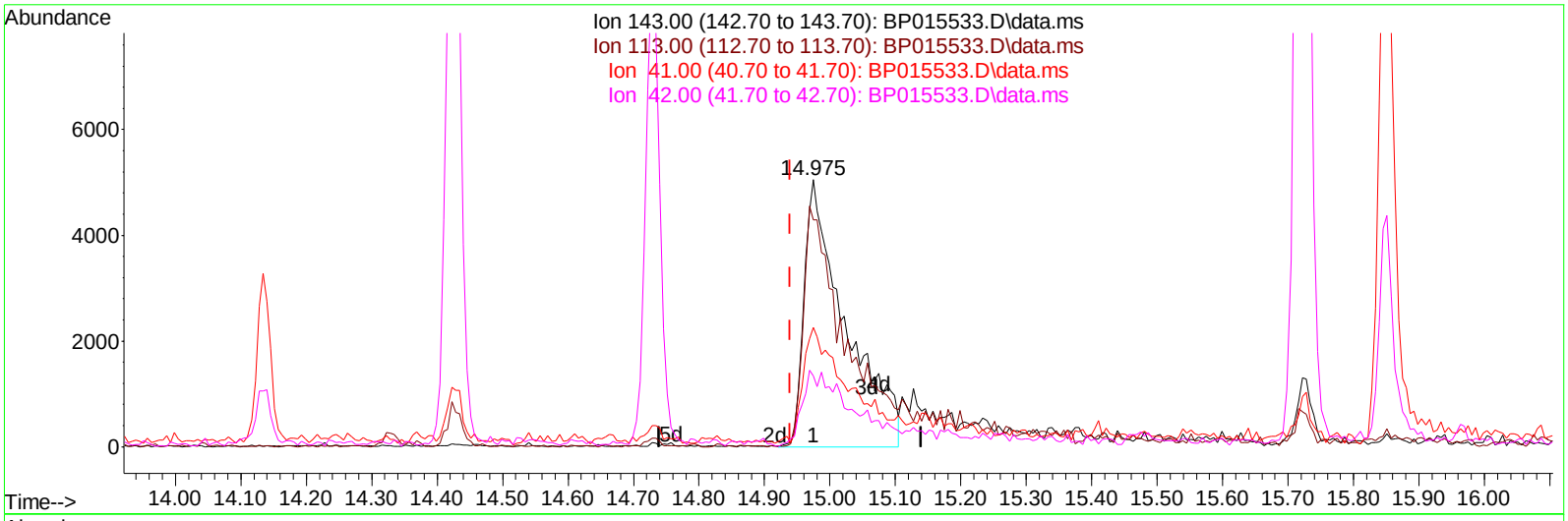

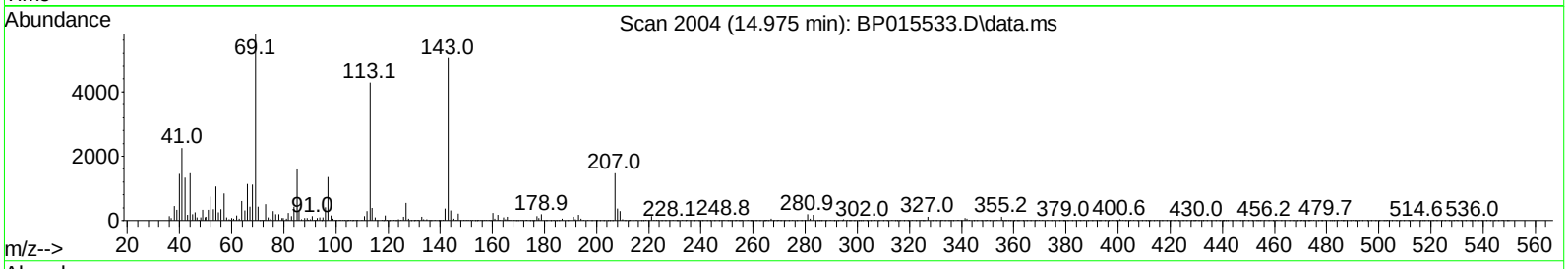

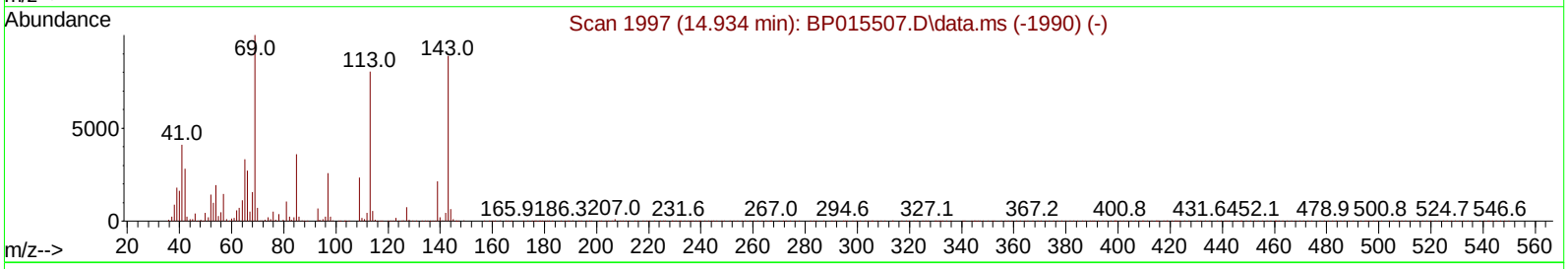

TIC: BP015533.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.035) 4.50 ng/ul m

response 22172

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	85.01
41.00	42.20	44.70
42.00	29.40	26.48

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	104251	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	499025	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	349189	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.487	188	836191	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	839724	20.000	ng/ul	0.00
88) Perylene-d12	24.116	264	898740	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	14696	5.220	ng/uL	0.00
4) Pyridine-d5	3.858	84	73221	10.013	ng/ul	0.01
7) Phenol-d5	7.246	99	83242	8.667	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	247313	43.117	ng/ul	0.00
11) 2-Chlorophenol-d4	7.611	132	237703	34.136	ng/ul	-0.01
15) 4-Methylphenol-d8	8.799	113	168764	21.410	ng/ul	-0.01
21) Nitrobenzene-d5	9.263	128	162644	41.514	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	181191	40.497	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	288805	35.386	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	329876	28.158	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1235976	44.886	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1314823	43.669	ng/ul	0.00
54) 4-Nitrophenol-d4	14.975	143	22172m	4.496	ng/ul	0.04
60) Fluorene-d10	15.722	176	1051816	45.298	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	199672	37.693	ng/ul	0.00
73) Anthracene-d10	17.587	188	1773200	46.607	ng/ul	0.00
81) Pyrene-d10	19.810	212	2197008	48.884	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2132876	47.297	ng/ul	0.00

Target Compounds Qvalue

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

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