Quantitation Report (Qedit)

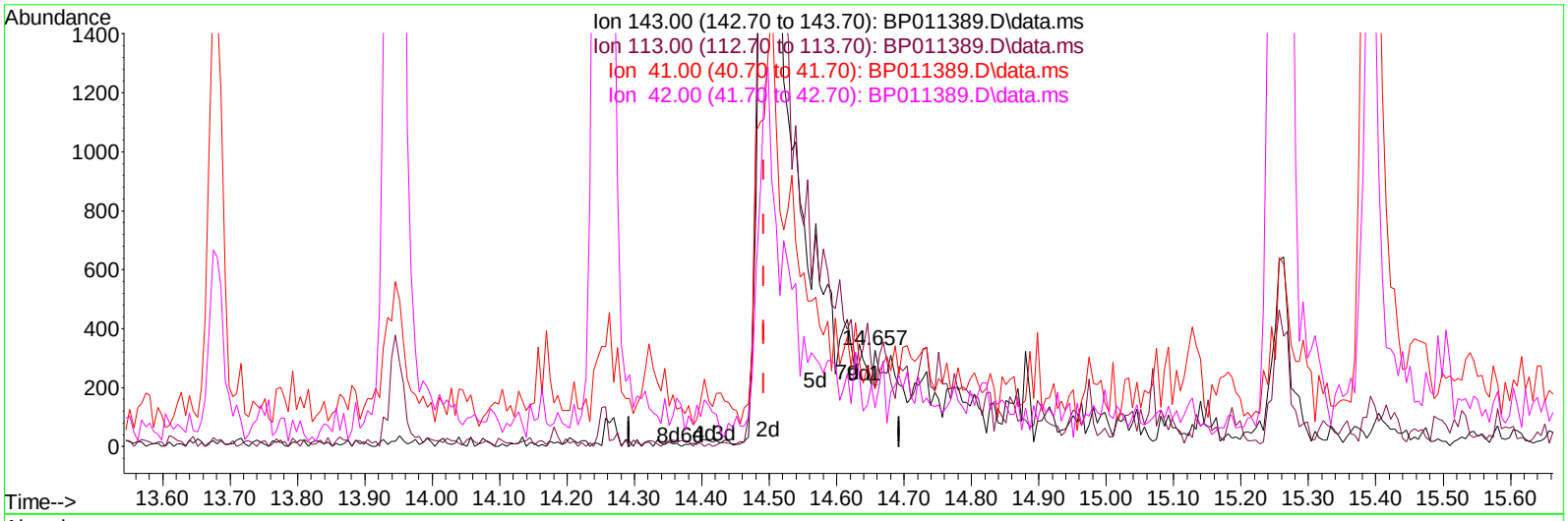
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011389.D
Acq On : 11 Aug 2022 16:11
Operator : CG/JU
Sample : N4128-17
Misc :
ALS Vial : 12 Sample Multiplier: 1

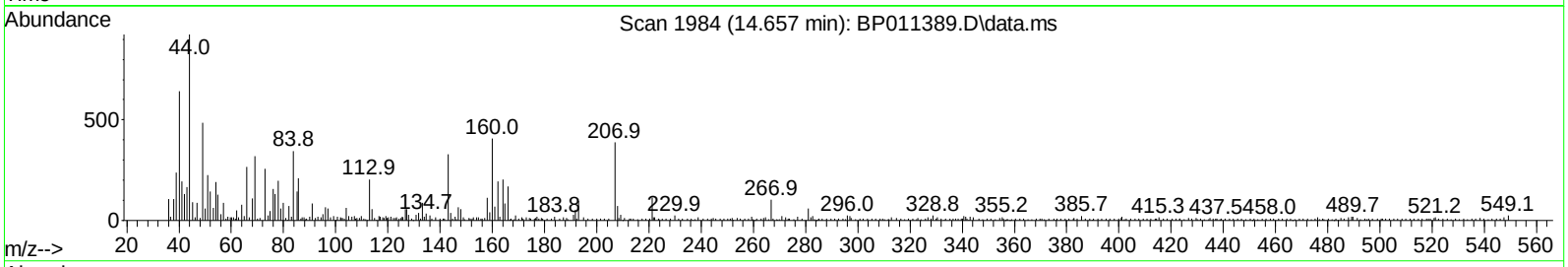
Quant Time: Aug 11 16:54:16 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

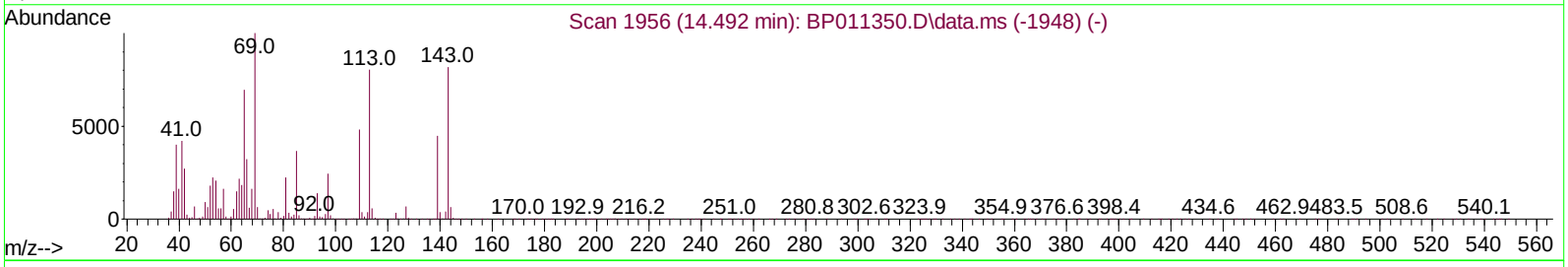
Instrument :
BNA_P
ClientSampleId :
BGBT3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011389.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.657min (+ 0.165) 0.02 ng/ul

response 56

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.20
41.00	54.40	59.45
42.00	35.50	40.24

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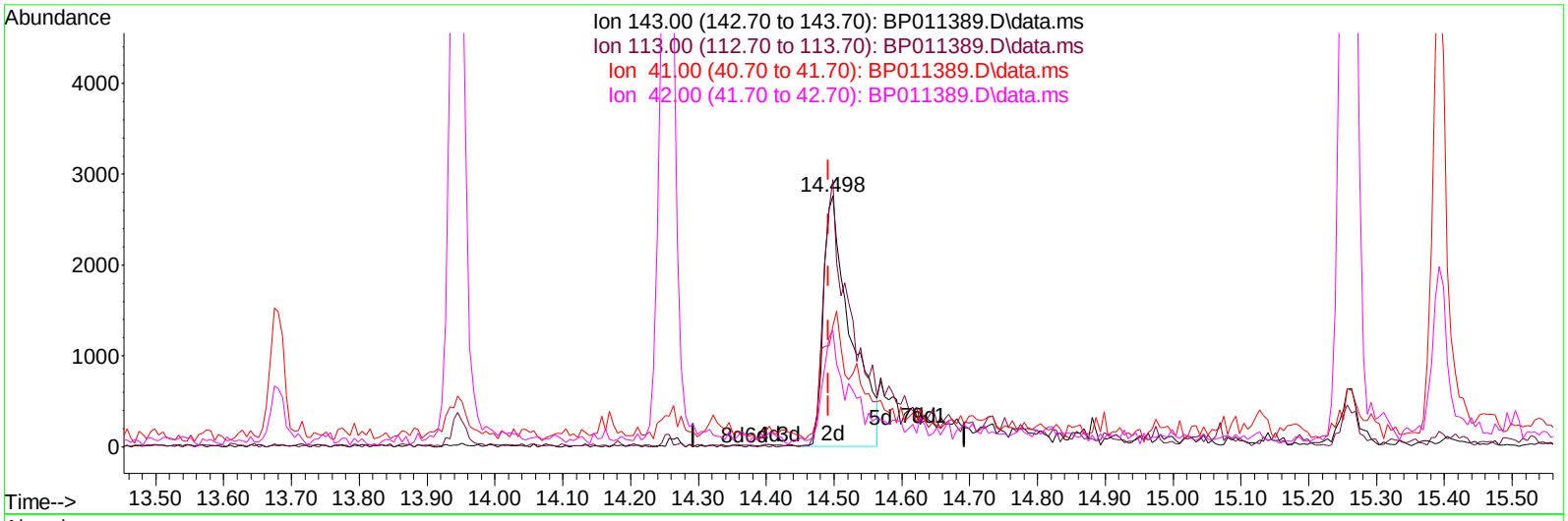

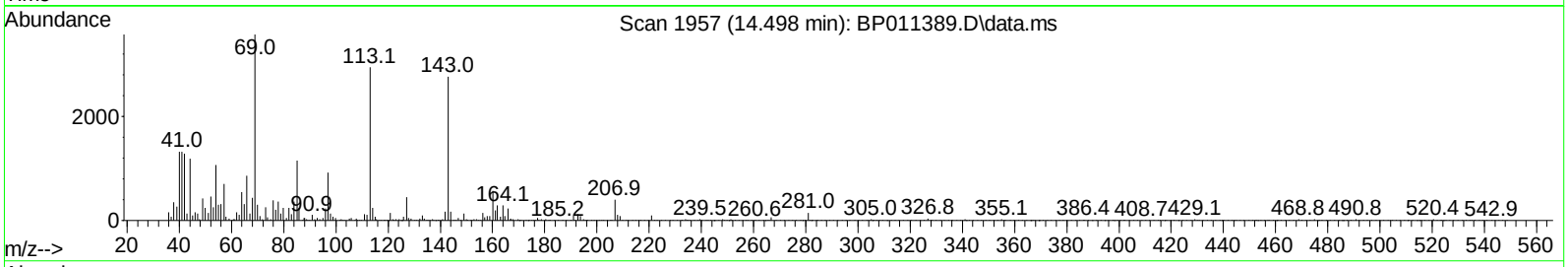

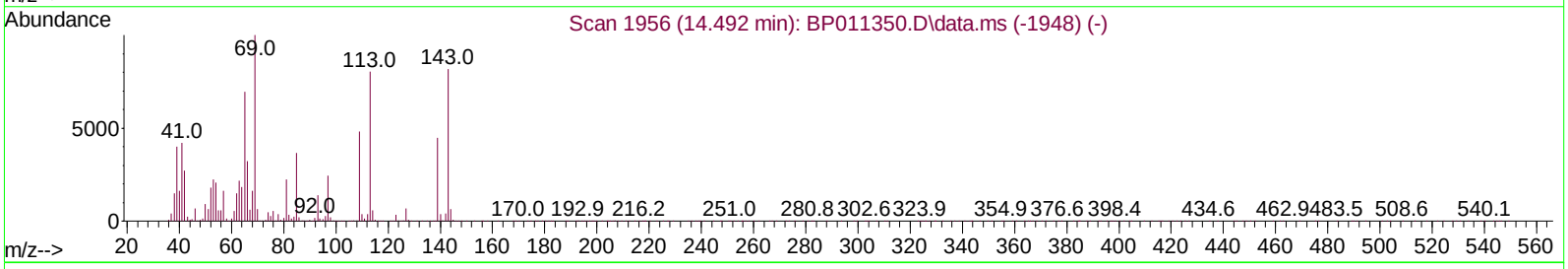

TIC: BP011389.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.73 ng/ul m

response 7572

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	106.57#
41.00	54.40	48.02
42.00	35.50	46.72#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	92446	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.369	136	386895	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	218966	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	452873	20.000	ng/ul	-0.01
79) Chrysene-d12	21.163	240	439386	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	459966	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	13245	4.392	ng/uL	0.00
4) Pyridine-d5	3.499	84	46216	6.076	ng/ul	0.00
7) Phenol-d5	6.763	99	46792	5.719	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	151454	26.109	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.110	132	139350	21.708	ng/ul	-0.01
15) 4-Methylphenol-d8	8.299	113	84838	13.207	ng/ul	-0.02
21) Nitrobenzene-d5	8.746	128	82137	31.194	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	77542	32.165	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.999	165	121566	24.971	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.522	131	173466	20.381	ng/ul	-0.01
46) Dimethylphthalate-d6	13.681	166	487933	32.116	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	552353	30.202	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	7572m	2.729	ng/ul	0.00
60) Fluorene-d10	15.257	176	409038	31.481	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.392	200	58673	27.893	ng/ul	0.00
73) Anthracene-d10	17.128	188	697303	34.256	ng/ul	-0.01
81) Pyrene-d10	19.392	212	820104	33.878	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	807160	34.797	ng/ul	0.00
Target Compounds						Qvalue

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

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