Quantitation Report (Qedit)

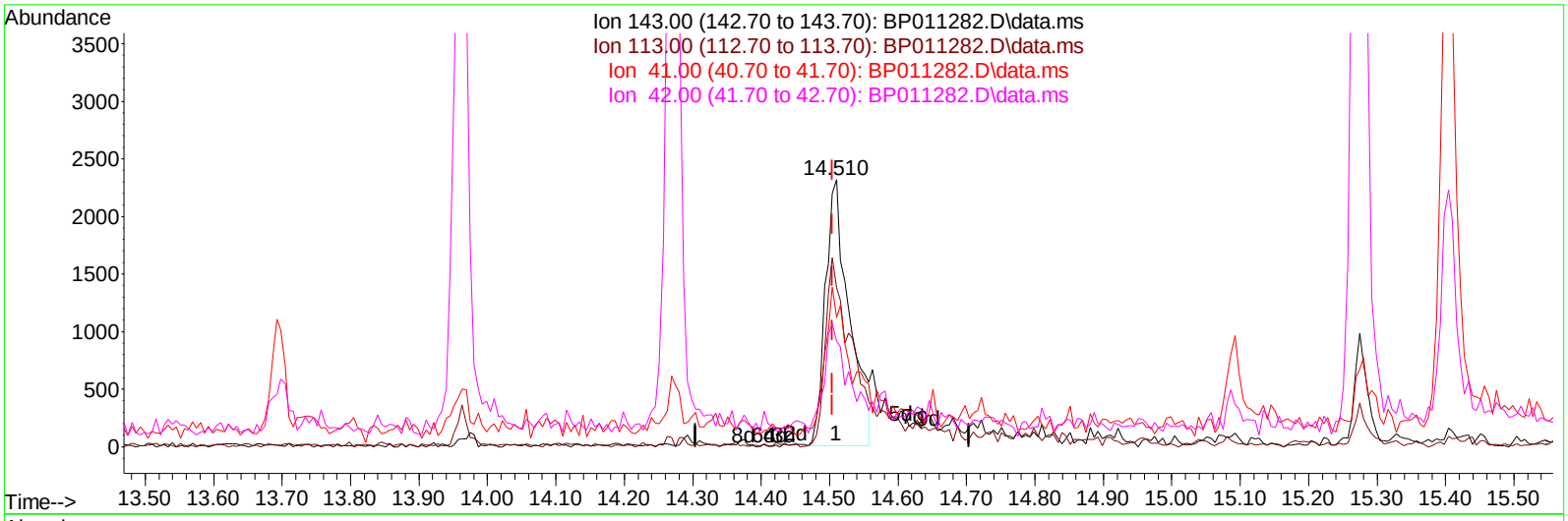
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011282.D
Acq On : 05 Aug 2022 17:12
Operator : CG/JU
Sample : N4026-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

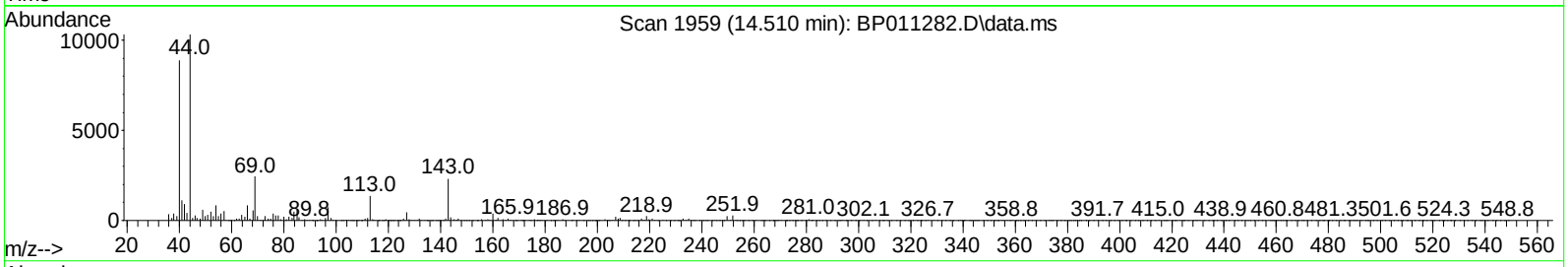
Quant Time: Aug 05 23:02:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 02 23:52:37 2022
Response via : Initial Calibration

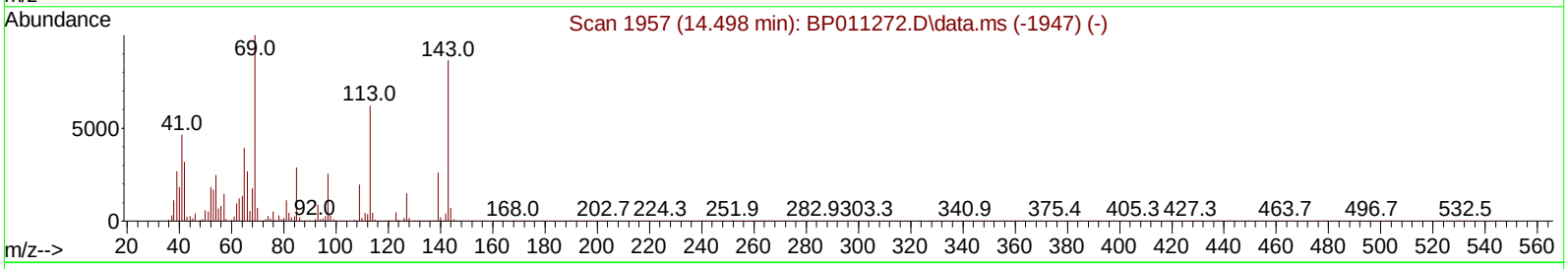
Instrument :
BNA_P
ClientSampleId :
GBJG9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/08/2022
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011282.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 2.42 ng/ul

response 5658

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	59.58#
41.00	54.40	48.36
42.00	35.50	39.39

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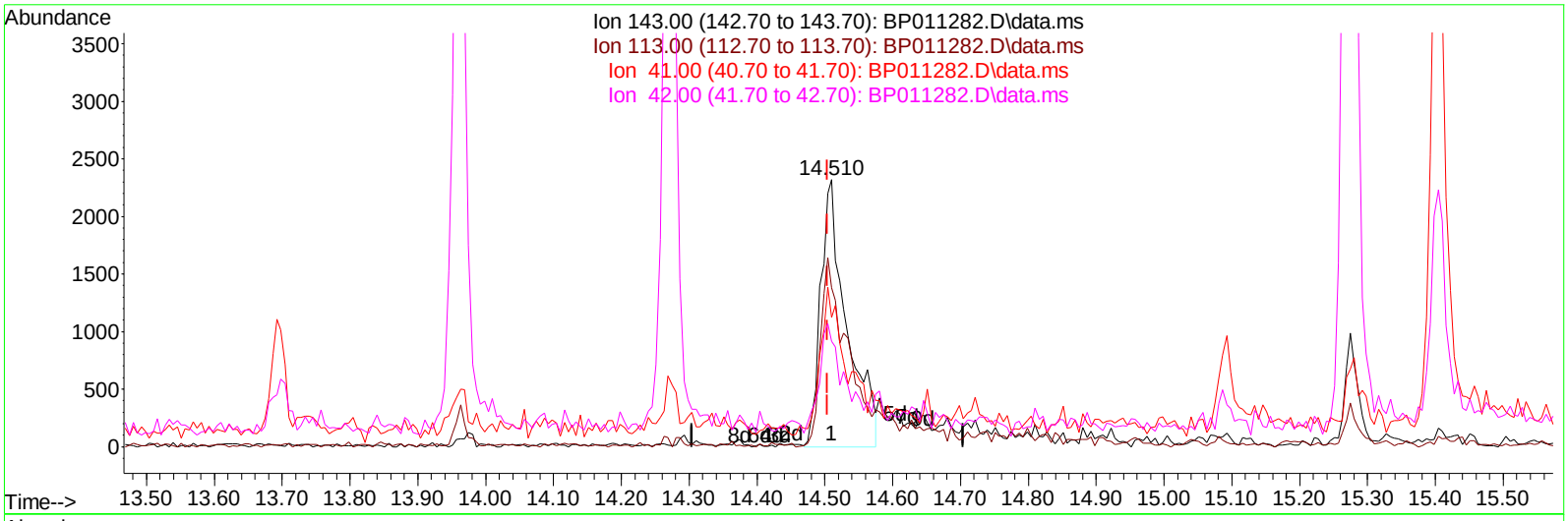

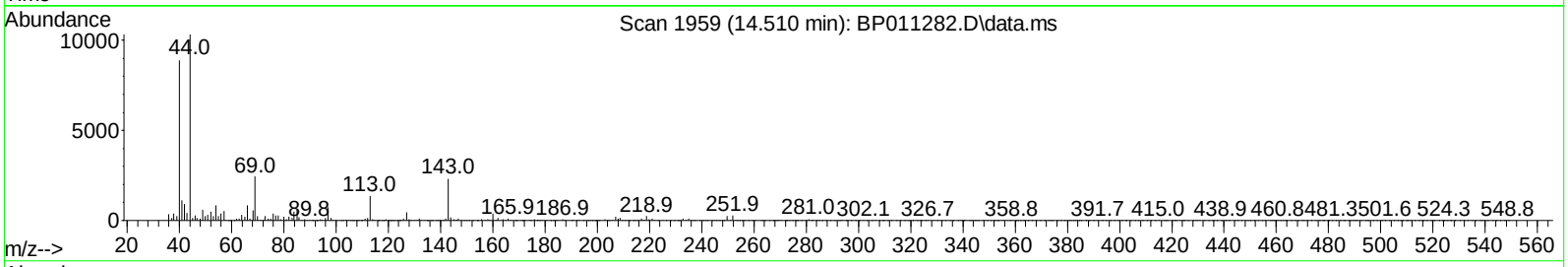

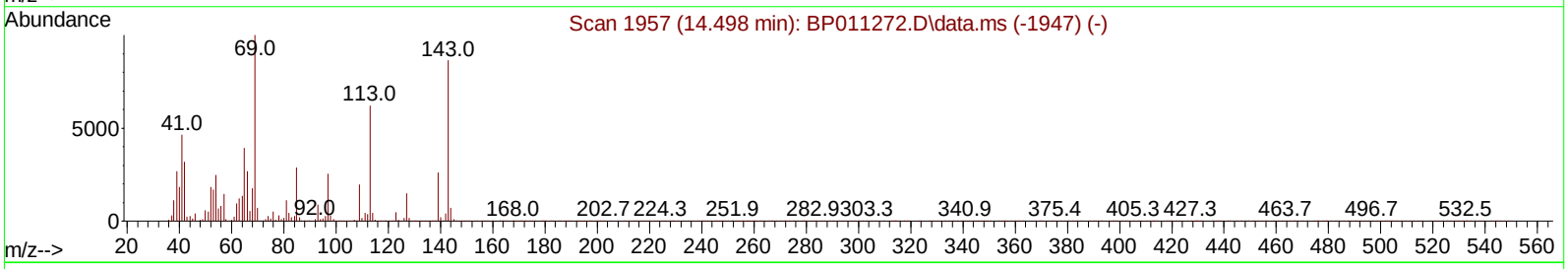

TIC: BP011282.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 2.66 ng/ul m

response 6207

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	59.58#
41.00	54.40	48.36
42.00	35.50	39.39

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	105305	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	398208	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.275	164	200473	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.045	188	372523	20.000	ng/ul	0.00
79) Chrysene-d12	21.168	240	551302	20.000	ng/ul	0.00
88) Perylene-d12	23.480	264	598981	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	3253	2.498	ng/uL	0.00
4) Pyridine-d5	3.511	84	22353	6.363	ng/ul	0.00
7) Phenol-d5	6.781	99	26416	3.010	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	178062	28.119	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	107026	14.962	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	49732	6.887	ng/ul	-0.01
21) Nitrobenzene-d5	8.763	128	120182	41.960	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	72639	27.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	100808	19.690	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	46008	5.809	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	330862	24.097	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	524076	30.337	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	6207m	2.658	ng/ul	0.00
60) Fluorene-d10	15.275	176	456403	39.065	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	53893	31.156	ng/ul	-0.01
73) Anthracene-d10	17.139	188	950612	58.063	ng/ul	0.00
81) Pyrene-d10	19.404	212	1134934	40.717	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	1258179	42.153	ng/ul	-0.01

Target Compounds Qvalue

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

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