Quantitation Report (Qedit)

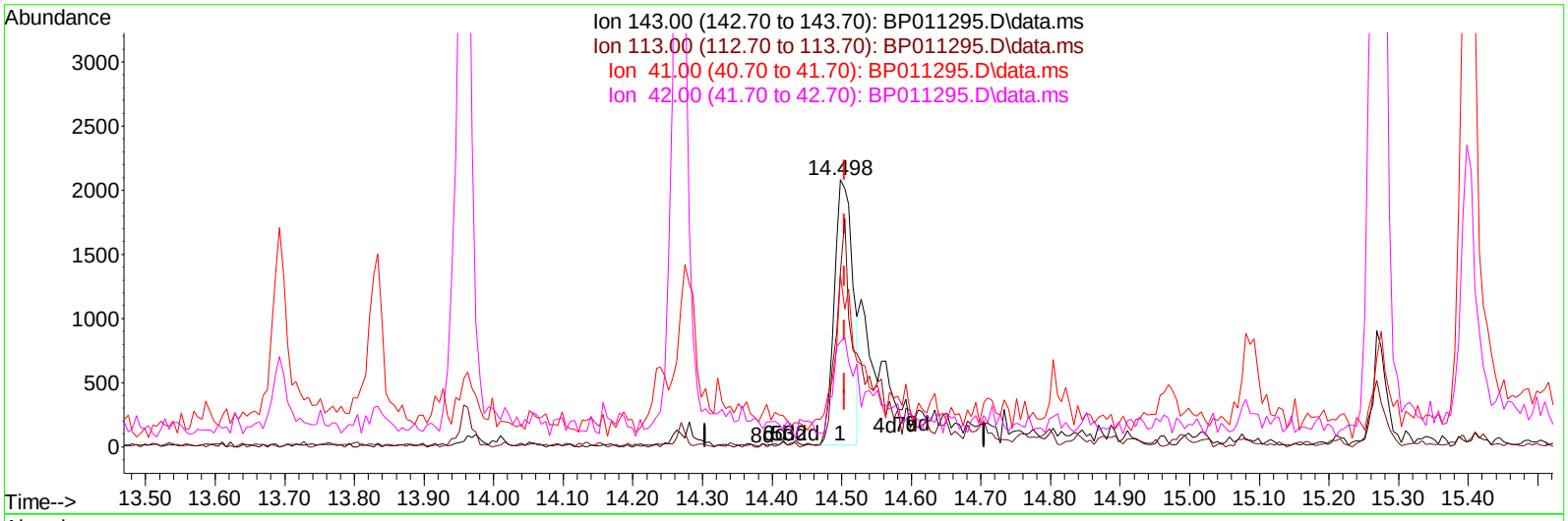
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011295.D
Acq On : 06 Aug 2022 00:45
Operator : CG/JU
Sample : N4026-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

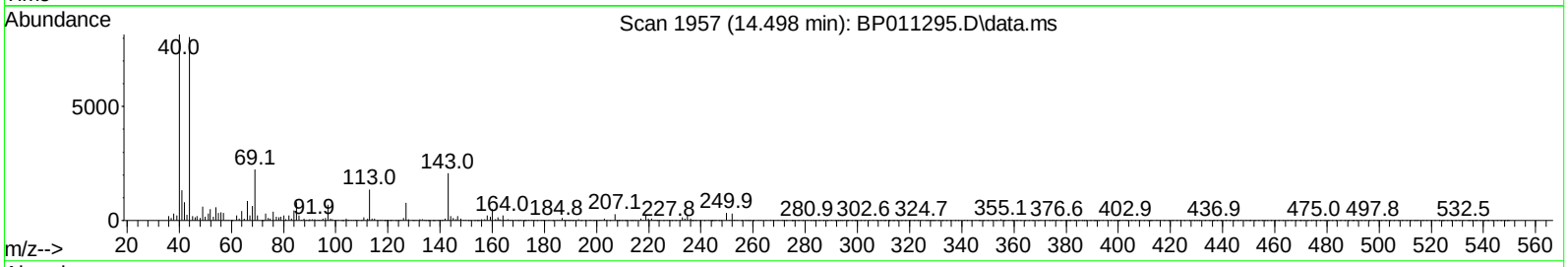
Quant Time: Aug 06 02:58:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 02 23:52:37 2022
Response via : Initial Calibration

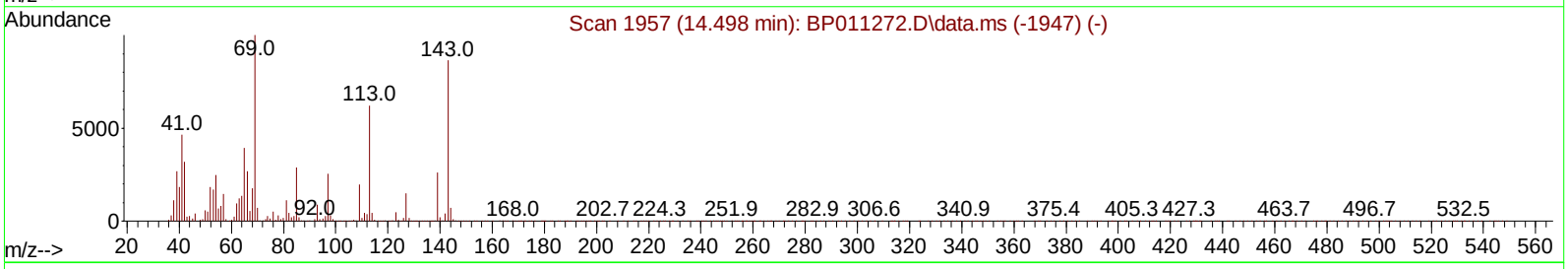
Instrument :
BNA_P
ClientSampleId :
GBJH0

Manual IntegrationsAPPROVED

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







TIC: BP011295.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 1.82 ng/ul

response 3873

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	65.45
41.00	54.40	64.30
42.00	35.50	39.21

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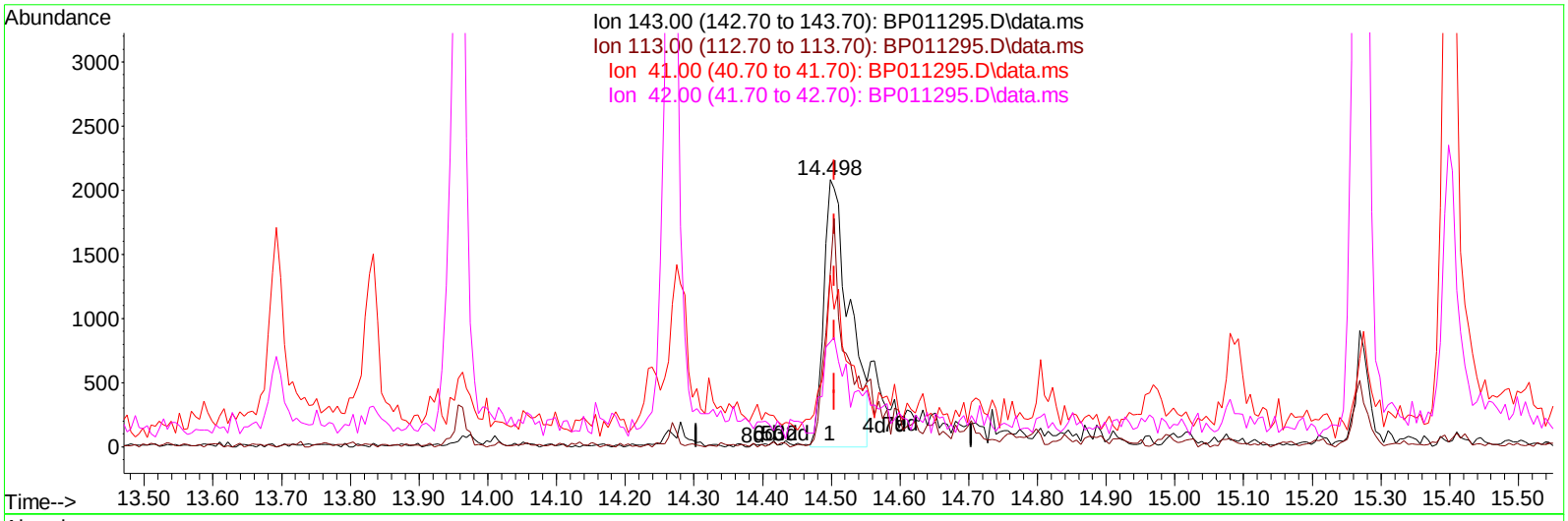

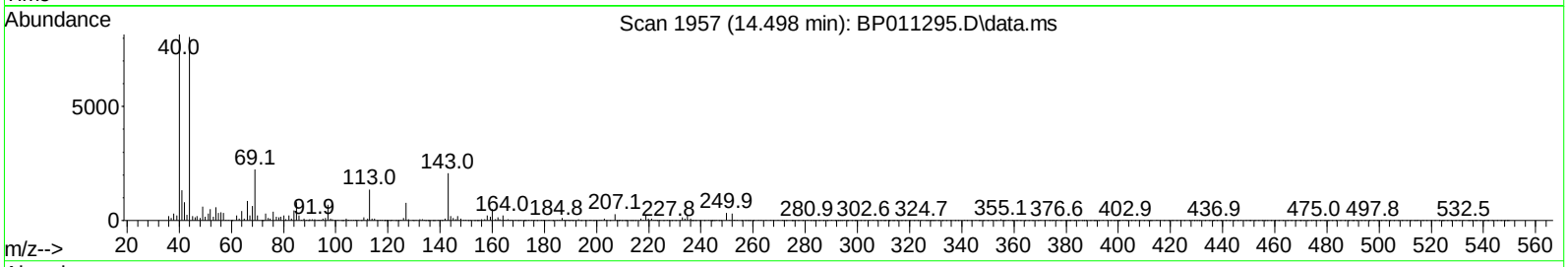

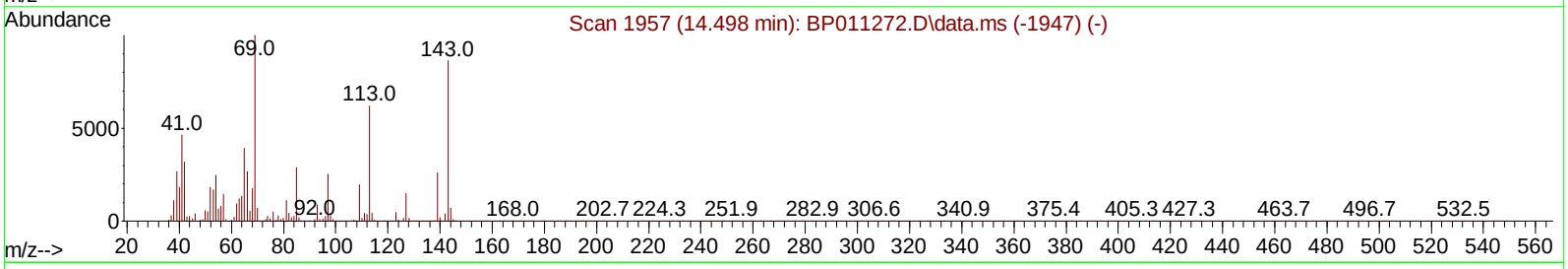

TIC: BP011295.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.51 ng/ul m

response 5331

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	65.45
41.00	54.40	64.30
42.00	35.50	39.21

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	88630	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	325362	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	182665	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188	376119	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	517329	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	534974	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2654	2.421	ng/uL	0.00
4) Pyridine-d5	3.511	84	11542	3.904	ng/ul	0.00
7) Phenol-d5	6.781	99	21149	2.863	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	147575	27.689	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	90916	15.101	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	42800	7.042	ng/ul	-0.01
21) Nitrobenzene-d5	8.757	128	110695	47.300	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	67577	30.806	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.016	165	88594	21.179	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	38469	5.944	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	329623	26.347	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	526597	33.454	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	5331m	2.506	ng/ul	0.00
60) Fluorene-d10	15.269	176	471191	44.263	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.398	200	58009	33.215	ng/ul	-0.02
73) Anthracene-d10	17.133	188	984377	59.550	ng/ul	-0.01
81) Pyrene-d10	19.398	212	1133760	43.345	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1116983	41.900	ng/ul	-0.02

Target Compounds Qvalue

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

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