Quantitation Report (Qedit)

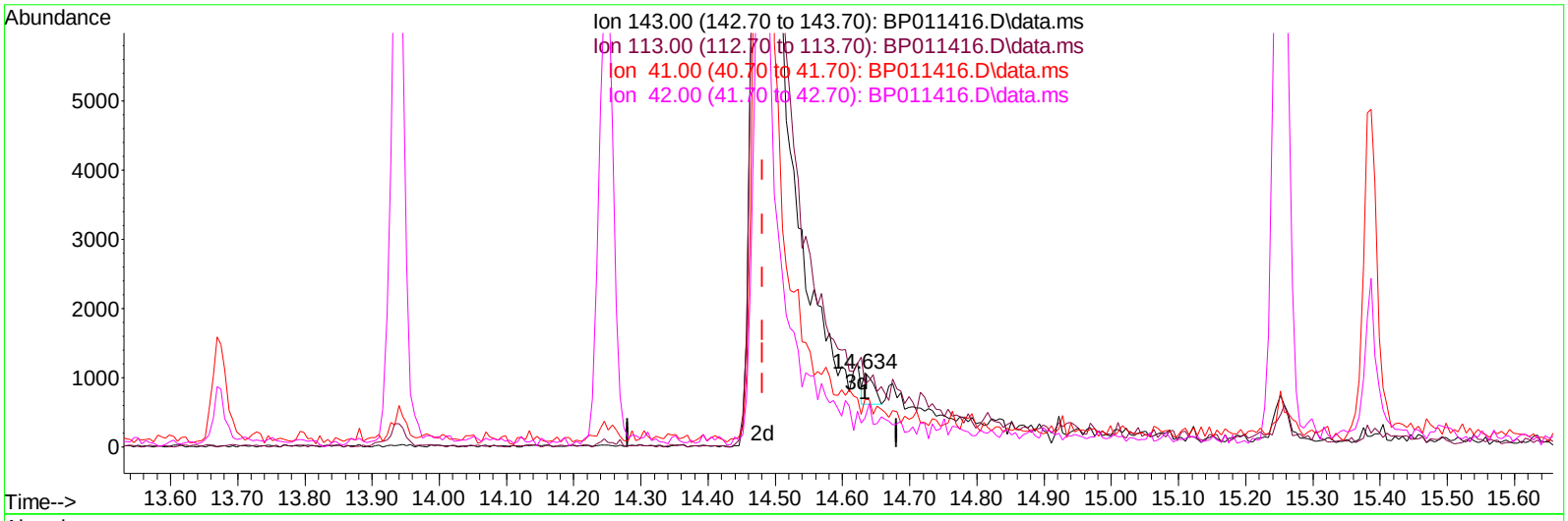
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011416.D
Acq On : 12 Aug 2022 08:50
Operator : CG/JU
Sample : PB146919BL
Misc :
ALS Vial : 40 Sample Multiplier: 1

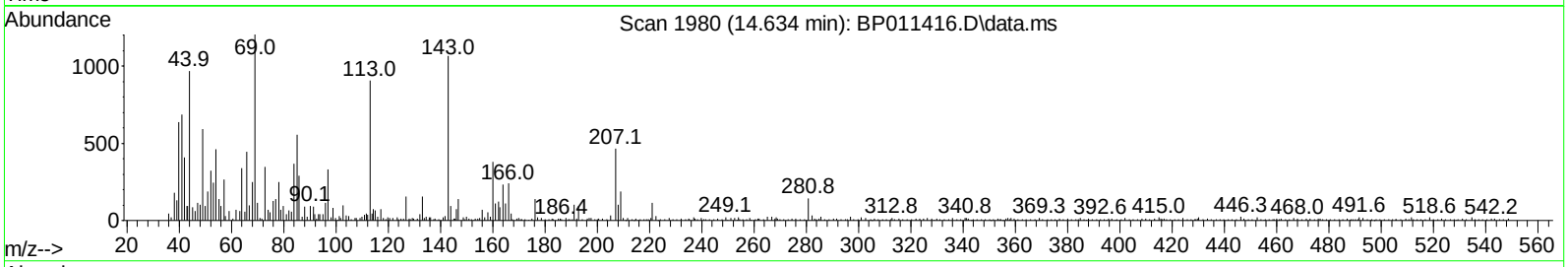
Quant Time: Aug 12 14:19:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

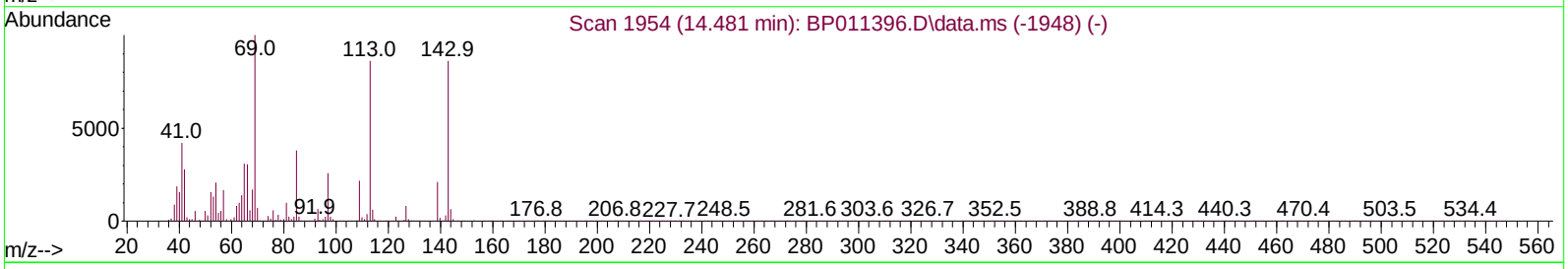
Instrument :
BNA_P
ClientSampleId :
SBLK919

Manual IntegrationsAPPROVED

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







TIC: BP011416.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.153) 0.18 ng/ul

response 457

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	84.90
41.00	54.40	64.26
42.00	35.50	38.46

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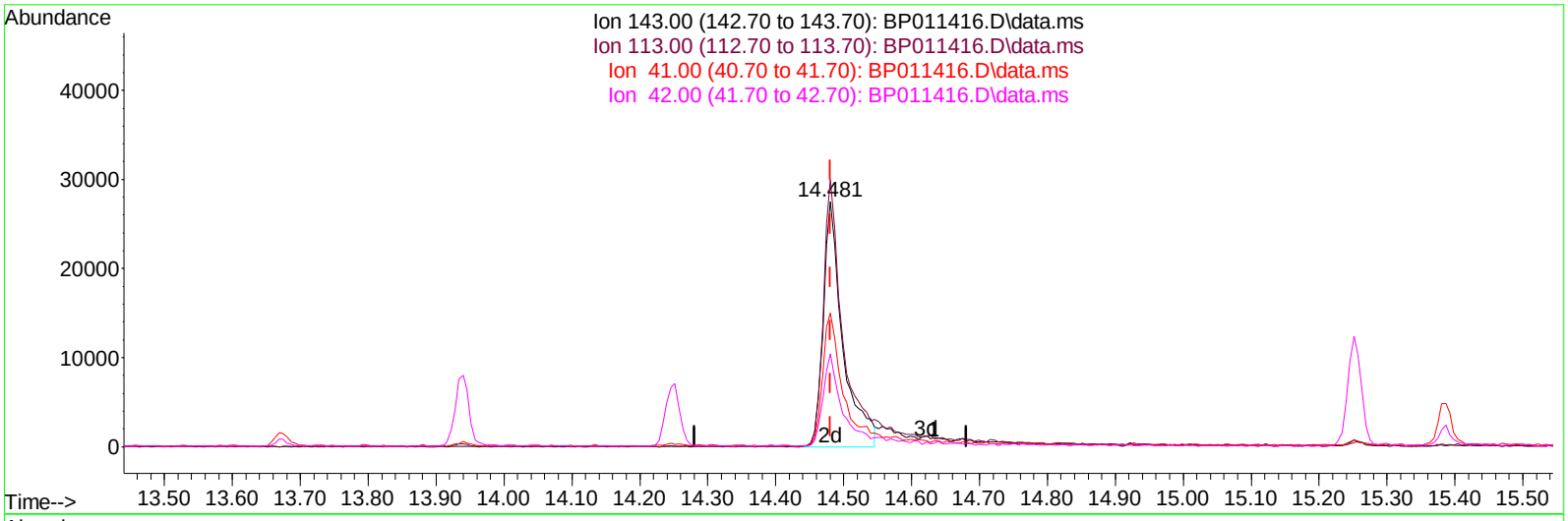

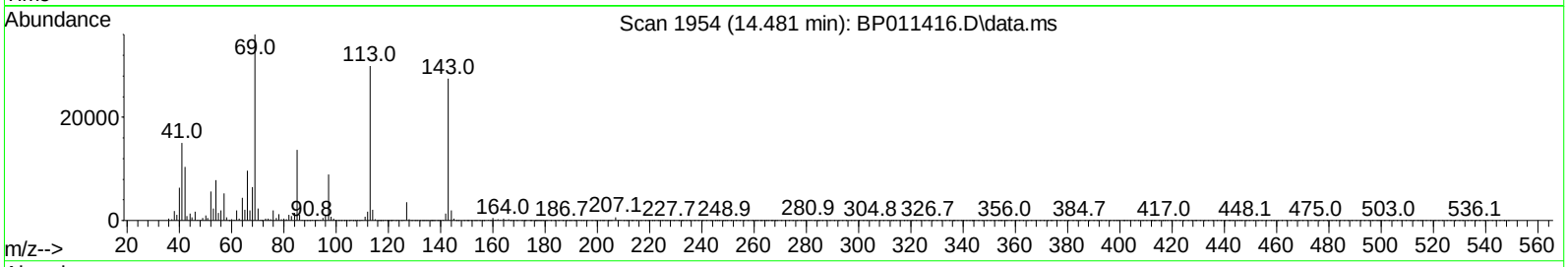

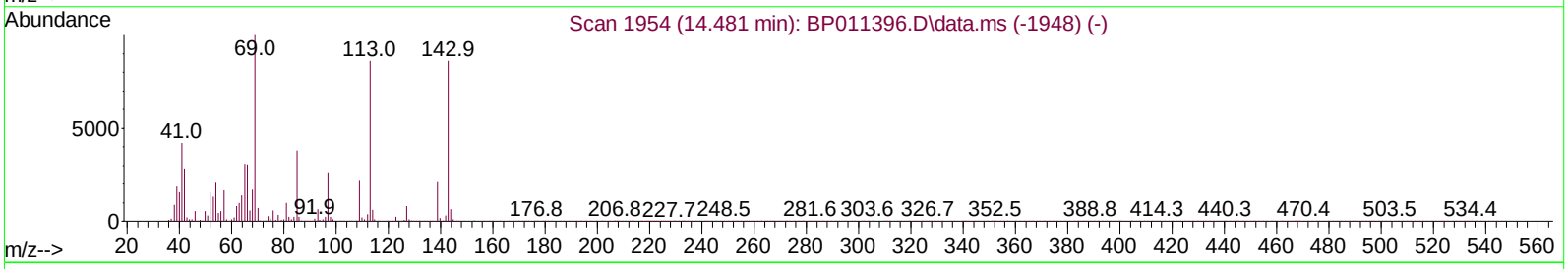

TIC: BP011416.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.000) 22.08 ng/ul m

response 54809

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	108.83#
41.00	54.40	54.64
42.00	35.50	37.96

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	79966	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	343248	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	195875	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	408283	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	391655	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	417818	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	15470	5.931	ng/uL	0.00
4) Pyridine-d5	3.487	84	195169	29.661	ng/ul	0.00
7) Phenol-d5	6.758	99	210474	29.741	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	148030	29.501	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	177094	31.894	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	165504	29.785	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	81248	34.780	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	80015	37.412	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	133976	31.019	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	200670	26.575	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	477057	35.102	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	539964	33.005	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	54809m	22.080	ng/ul	0.00
60) Fluorene-d10	15.251	176	380589	32.745	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	53816	28.378	ng/ul	0.00
73) Anthracene-d10	17.122	188	615242	33.526	ng/ul	0.00
81) Pyrene-d10	19.386	212	716125	33.188	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	700291	33.235	ng/ul	0.00

Target Compounds Qvalue

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

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