Quantitation Report (Qedit)

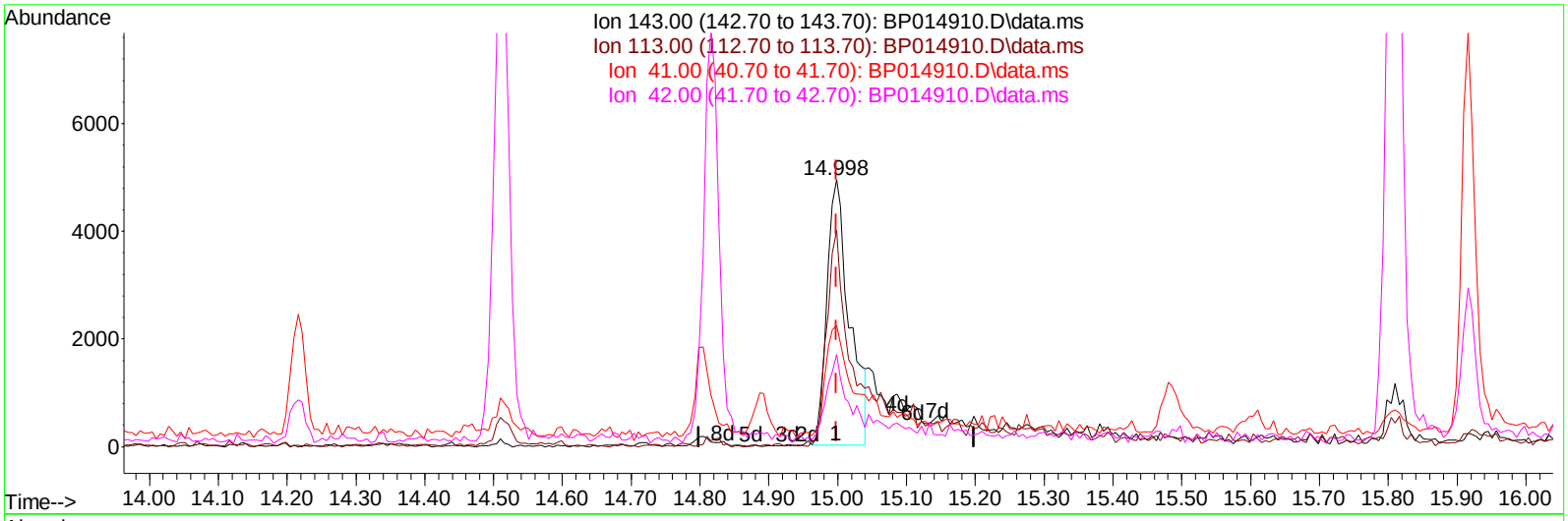
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050123\
Data File : BP014910.D
Acq On : 01 May 2023 12:26
Operator : CG/JU
Sample : 02551-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

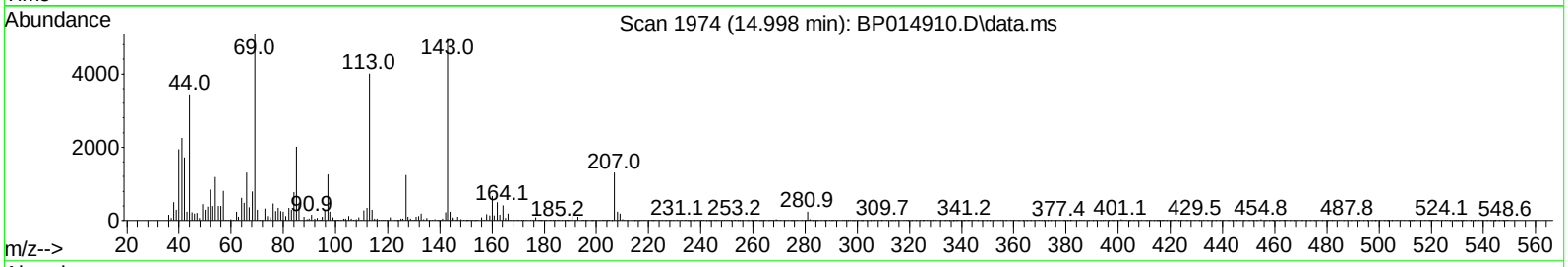
Quant Time: May 02 00:19:11 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 29 00:40:33 2023
Response via : Initial Calibration

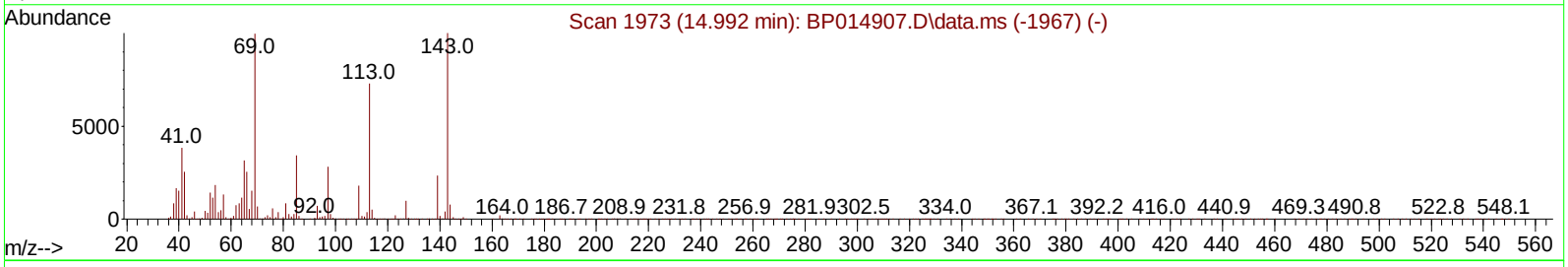
Instrument :
BNA_P
ClientSampleId :
BHE78

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/02/2023
Supervised By :Jagrut Upadhyay 05/02/2023







TIC: BP014910.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 2.57 ng/ul

response 11170

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	80.92
41.00	36.10	45.47#
42.00	26.80	34.60#

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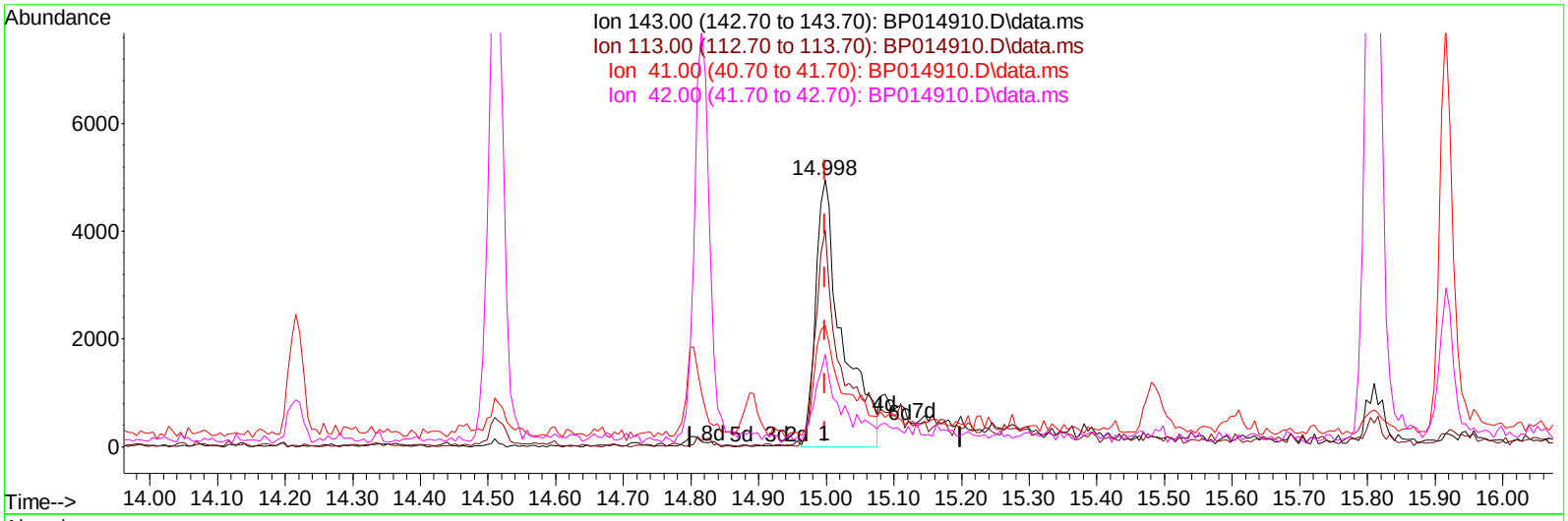

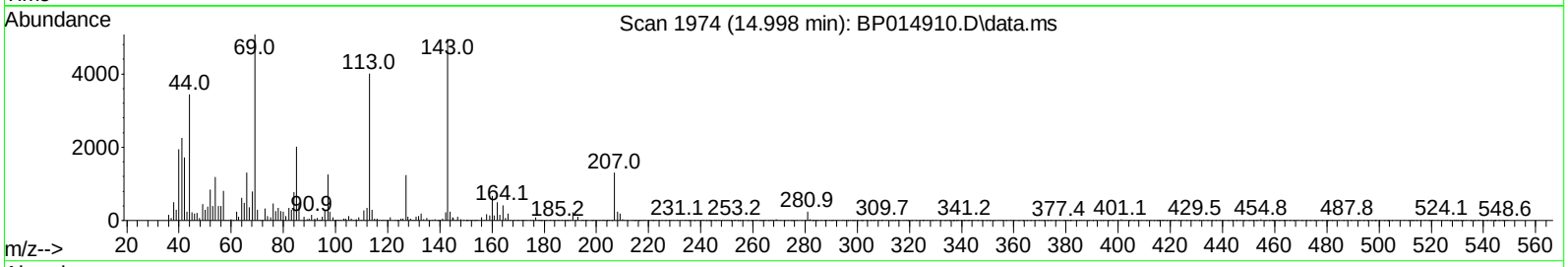

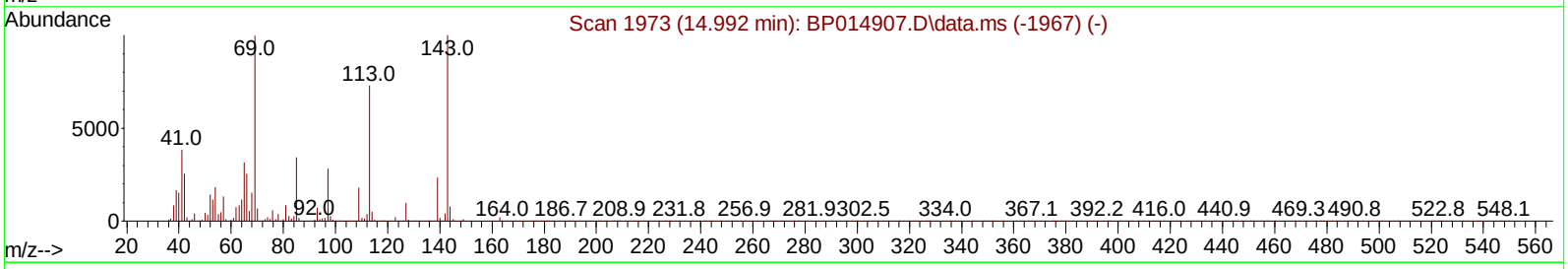

TIC: BP014910.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 3.10 ng/ul m

response 13483

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	80.92
41.00	36.10	45.47#
42.00	26.80	34.60#

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050123\
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Operator : CG/JU
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Misc :
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Quant Time: May 02 00:31:24 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 29 00:40:33 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.175	152	171873	20.000	ng/ul	0.00
20) Naphthalene-d8	11.010	136	712834	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.816	164	407474	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.575	188	814239	20.000	ng/ul	0.00
79) Chrysene-d12	21.657	240	691073	20.000	ng/ul	0.00
88) Perylene-d12	24.262	264	796987	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.464	96	24922	5.226	ng/uL	0.00
4) Pyridine-d5	3.911	84	78412	6.840	ng/ul	0.00
7) Phenol-d5	7.310	99	108729	8.083	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.487	67	311549	36.926	ng/ul	0.00
11) 2-Chlorophenol-d4	7.693	132	305826	30.026	ng/ul	0.00
15) 4-Methylphenol-d8	8.875	113	190652	18.600	ng/ul	0.00
21) Nitrobenzene-d5	9.346	128	172126	33.630	ng/ul	0.00
24) 2-Nitrophenol-d4	10.075	143	166717	35.484	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.616	165	304578	30.184	ng/ul	0.00
31) 4-Chloroaniline-d4	11.140	131	404157	27.702	ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	1165781	39.921	ng/ul	0.00
49) Acenaphthylene-d8	14.510	160	1310733	37.576	ng/ul	0.00
54) 4-Nitrophenol-d4	14.998	143	13483m	3.103	ng/ul	0.00
60) Fluorene-d10	15.810	176	1001094	39.328	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.916	200	122359	35.527	ng/ul	0.00
73) Anthracene-d10	17.675	188	1702497	45.783	ng/ul	0.00
81) Pyrene-d10	19.892	212	1957333	50.995	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.092	264	1831412	46.524	ng/ul	0.00

Target Compounds Qvalue

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

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