Quantitation Report (Qedit)

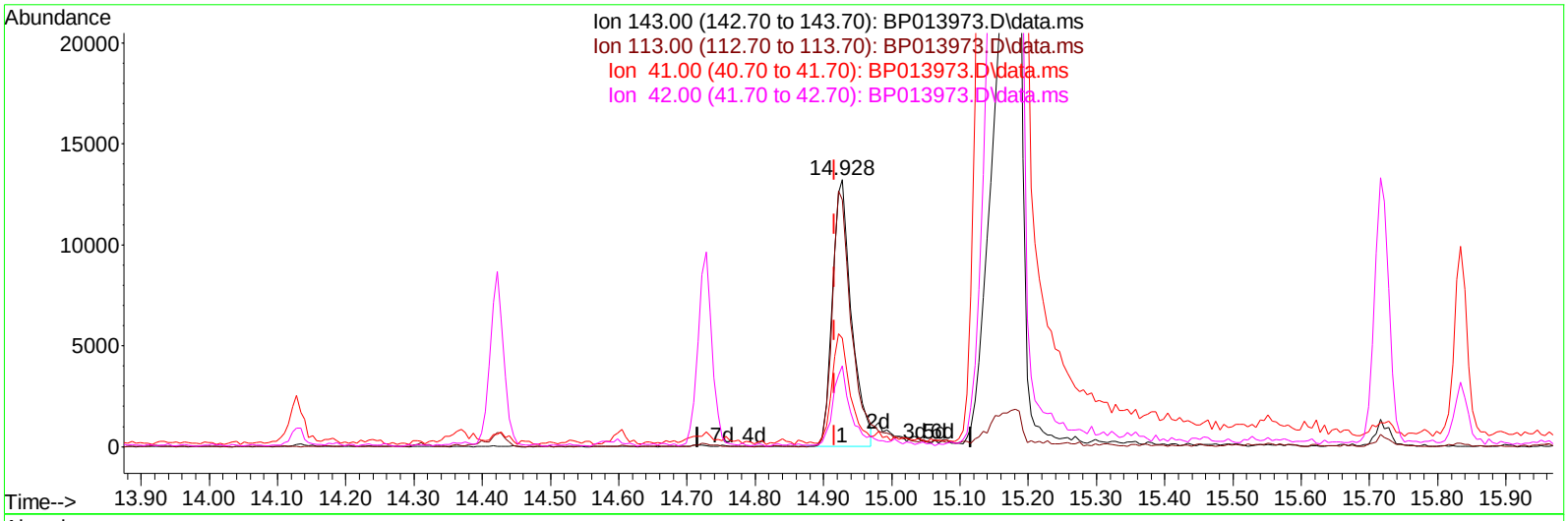
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013973.D
Acq On : 08 Mar 2023 12:37
Operator : CG/JU
Sample : 01729-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

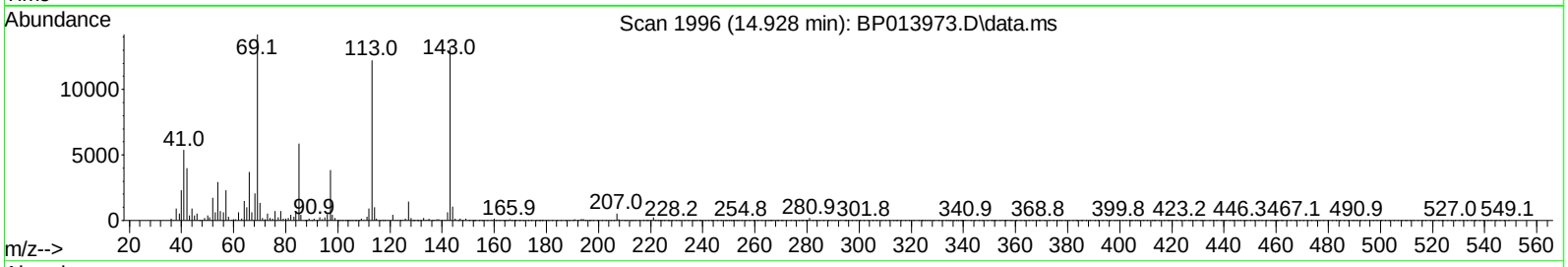
Quant Time: Mar 09 01:12:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 09 01:10:12 2023
Response via : Initial Calibration

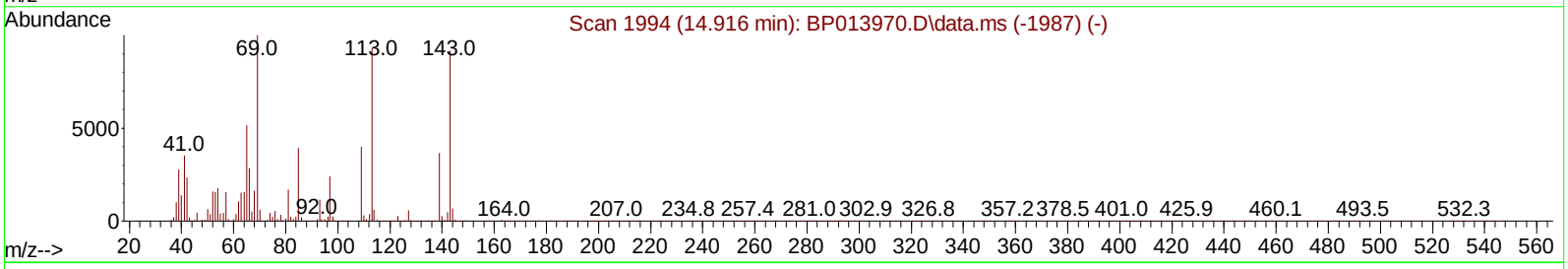
Instrument :
BNA_P
ClientSampleId :
C0BY9

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/09/2023
Supervised By :Jagrut Upadhyay 03/09/2023







TIC: BP013973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.012) 4.02 ng/ul

response 25682

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	92.45
41.00	40.20	40.63
42.00	28.50	30.38

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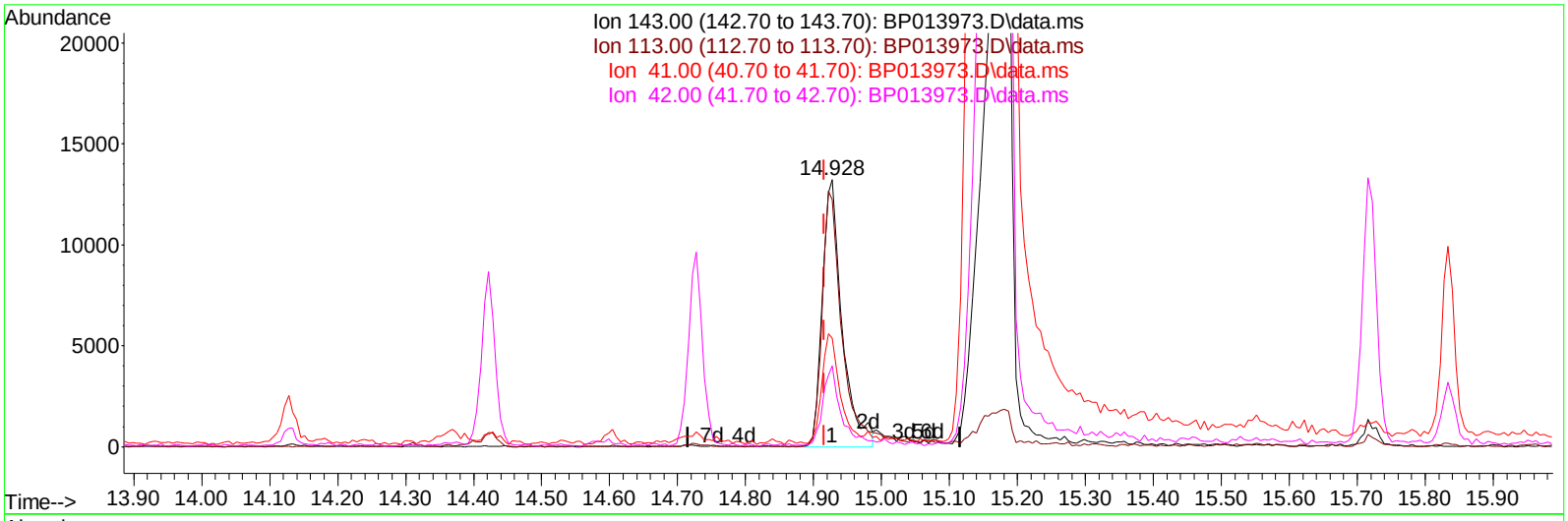

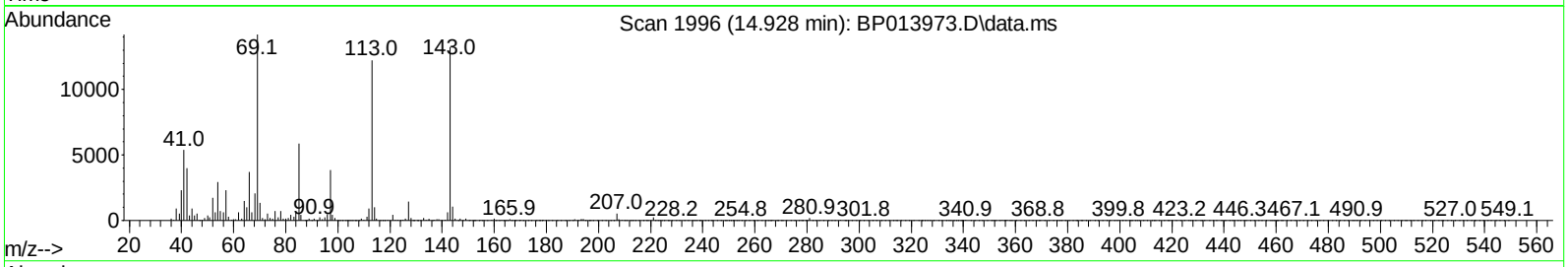

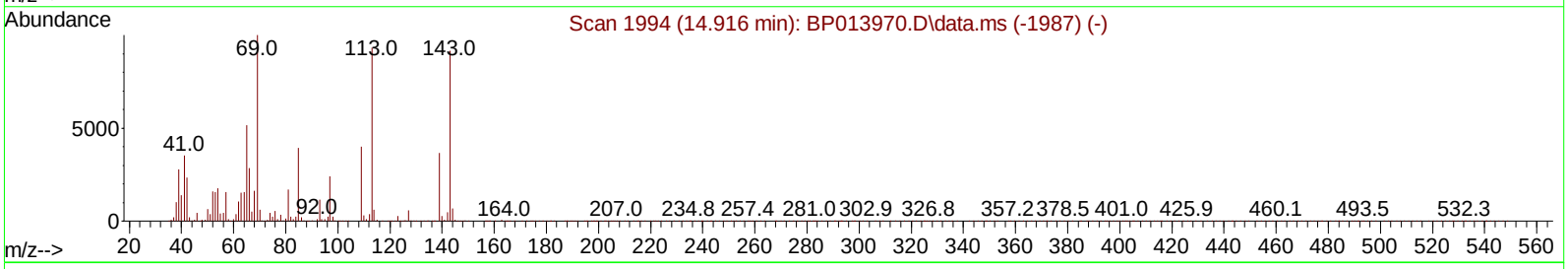

TIC: BP013973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.012) 4.19 ng/ul m

response 26744

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	92.45
41.00	40.20	40.63
42.00	28.50	30.38

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	138120	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	582769	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	426913	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.481	188	1011751	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	1141677	20.000	ng/ul	0.00
88) Perylene-d12	24.104	264	1210099	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.440	96	14516	3.971	ng/uL	0.00
4) Pyridine-d5	3.881	84	69982	6.914	ng/ul	0.01
7) Phenol-d5	7.240	99	64567	5.297	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	163061	22.939	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	174686	18.811	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	132208	13.279	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	106492	22.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.987	143	121482	21.825	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	256727	24.513	ng/ul	0.00
31) 4-Chloroaniline-d4	11.052	131	265595	19.163	ng/ul	0.00
46) Dimethylphthalate-d6	14.128	166	946741	26.131	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	940693	24.509	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	26744m	4.189	ng/ul	0.01
60) Fluorene-d10	15.722	176	790539	25.746	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.834	200	175172	22.908	ng/ul	0.00
73) Anthracene-d10	17.581	188	1384545	28.189	ng/ul	0.00
81) Pyrene-d10	19.804	212	1719884	27.314	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.939	264	1784797	27.728	ng/ul	0.01

Target Compounds Qvalue

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

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