Quantitation Report (Qedit)

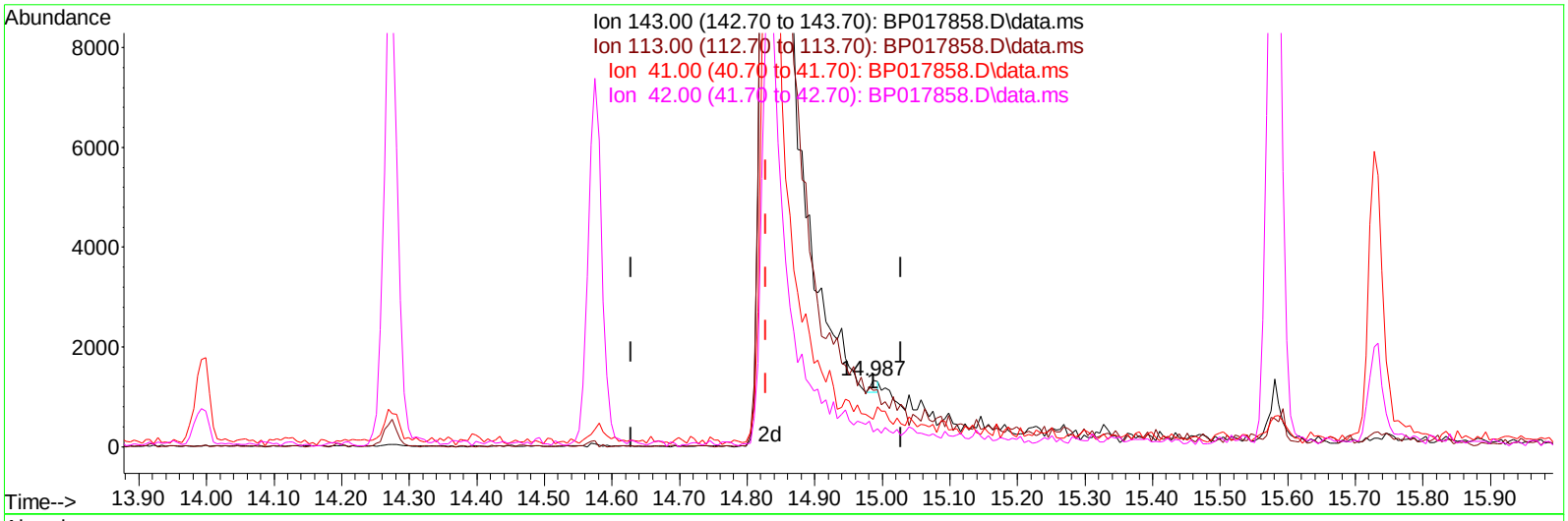
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017858.D
Acq On : 01 Nov 2023 20:03
Operator : MA/JU
Sample : PB156543BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

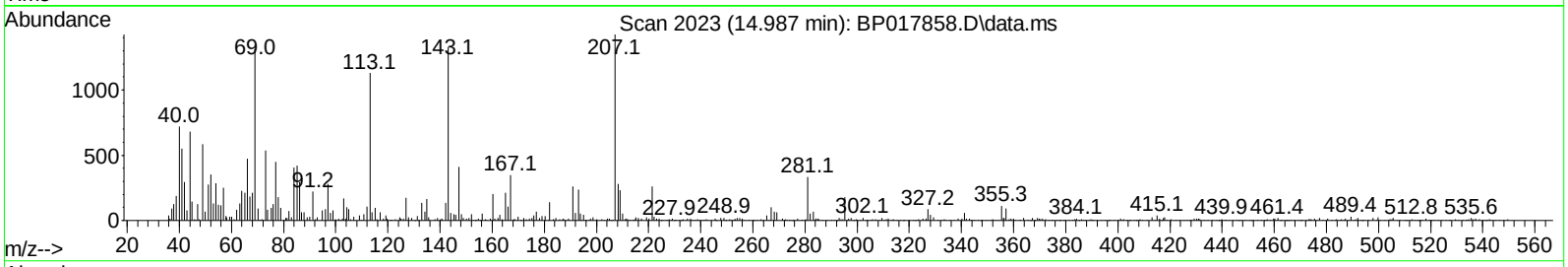
Quant Time: Nov 02 01:15:30 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 19 01:32:15 2023
Response via : Initial Calibration

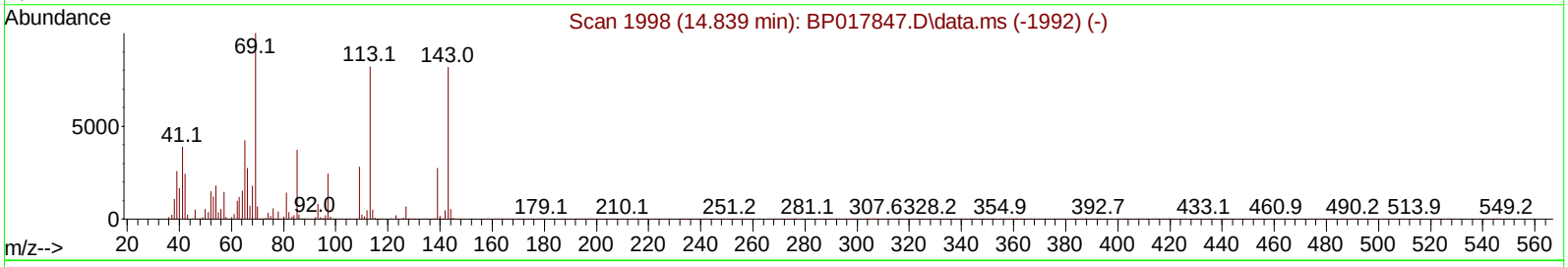
Instrument :
BNA_P
ClientSampleId :
SBLK543

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/03/2023
Supervised By :mohammad ahmed 11/04/2023







TIC: BP017858.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.987min (+ 0.159) 0.06 ng/ul

response 191

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	84.89
41.00	39.20	41.58
42.00	25.60	22.11

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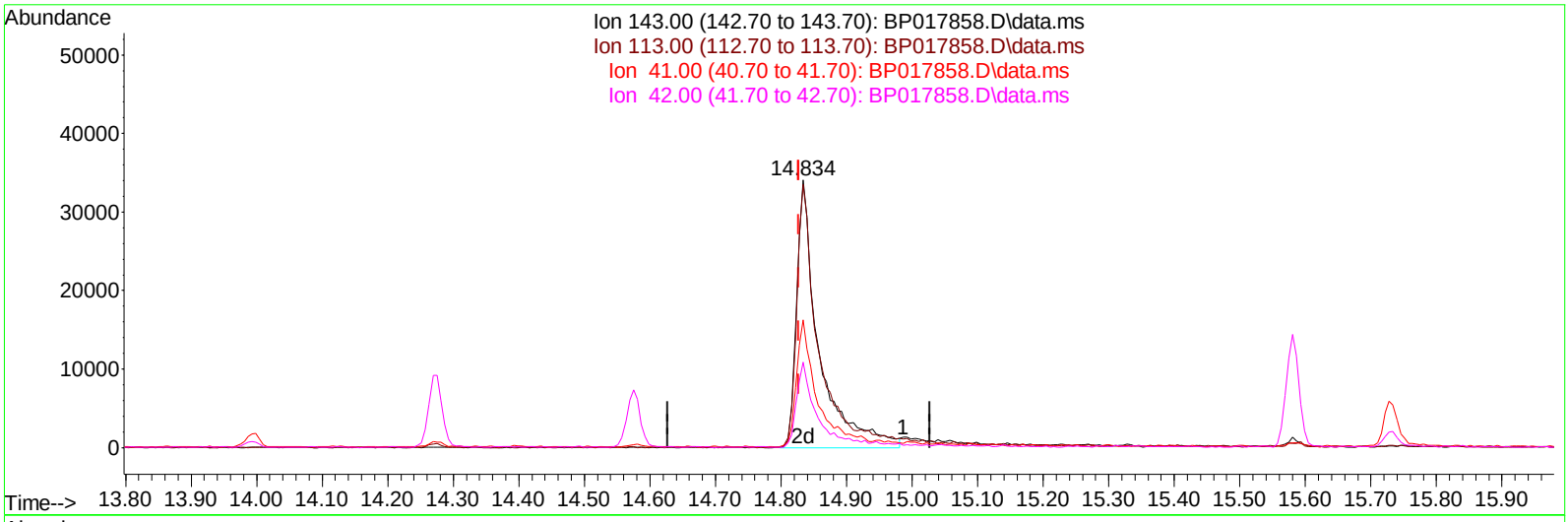

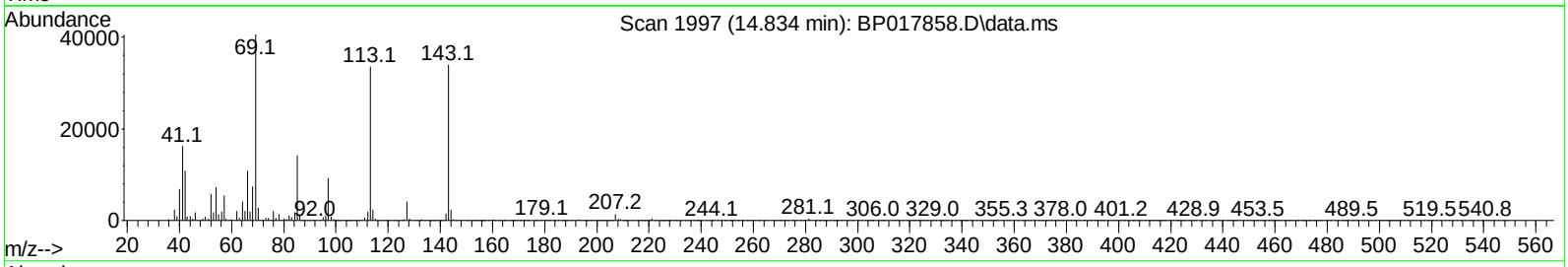

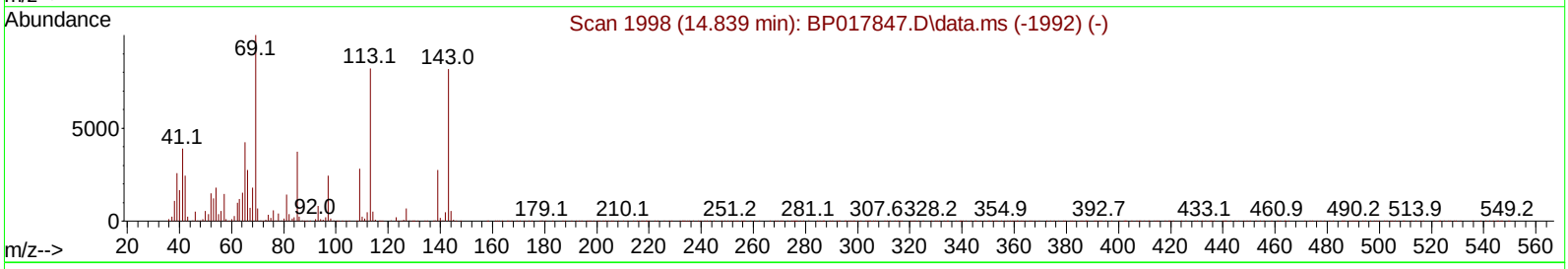

TIC: BP017858.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.006) 27.00 ng/ul m

response 81379

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	98.43#
41.00	39.20	47.68#
42.00	25.60	31.94#

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017858.D
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Misc :
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Quant Time: Nov 02 02:04:09 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	118952	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	447381	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.575	164	251955	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.369	188	511956	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	423280	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	444672	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	21325	6.669	ng/uL	0.02
4) Pyridine-d5	3.705	84	249837	29.843	ng/ul	0.01
7) Phenol-d5	7.040	99	338035	34.150	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	219681	35.936	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	282671	35.012	ng/ul	-0.01
15) 4-Methylphenol-d8	8.605	113	265552	34.162	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	129725	35.784	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.799	143	146375	36.544	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	245489	31.244	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.887	131	322439	31.358	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	734852	34.743	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	811185	34.390	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	81379m	27.003	ng/ul	0.00
60) Fluorene-d10	15.581	176	604703	33.074	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	96223	28.534	ng/ul	0.00
73) Anthracene-d10	17.469	188	938566	35.875	ng/ul	0.00
81) Pyrene-d10	19.733	212	1037338	39.089	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	901574	36.402	ng/ul	0.00

Target Compounds Qvalue

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

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