Quantitation Report (Qedit)

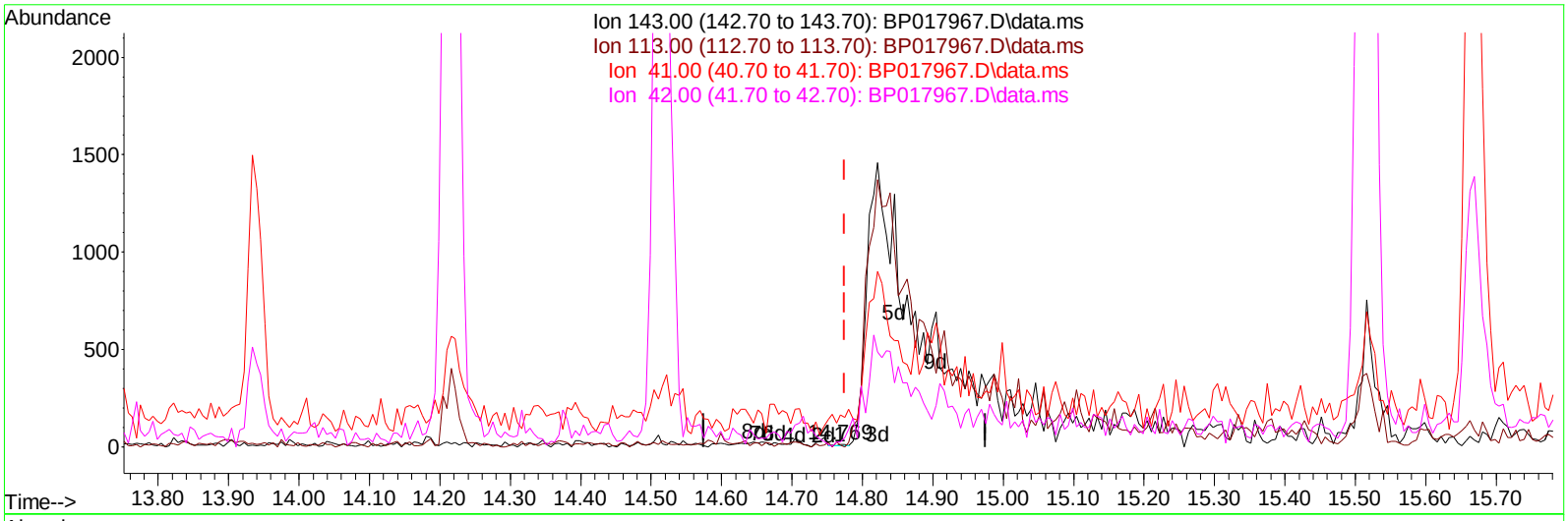
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
Data File : BP017967.D
Acq On : 08 Nov 2023 18:00
Operator : MA/JU
Sample : 05013-08
Misc :
ALS Vial : 47 Sample Multiplier: 1

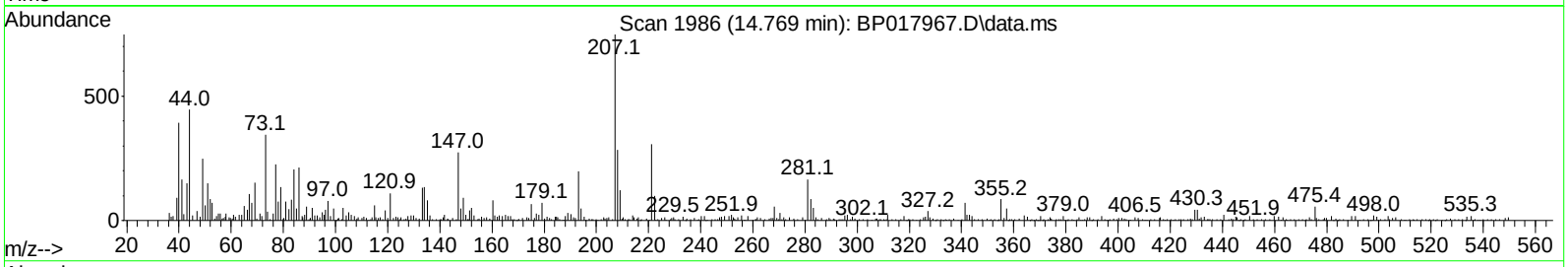
Quant Time: Nov 08 21:44:00 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 07 03:40:56 2023
Response via : Initial Calibration

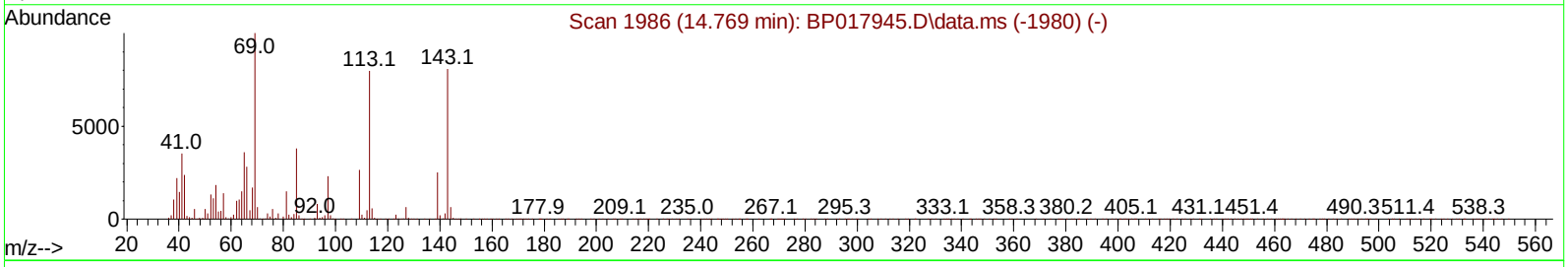
Instrument :
BNA_P
ClientSampleId :
A4995

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/09/2023
Supervised By :mohammad ahmed 11/11/2023







TIC: BP017967.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.006) 0.00 ng/ul

response 8

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	61.54#
41.00	43.00	1269.23#
42.00	28.20	192.31#

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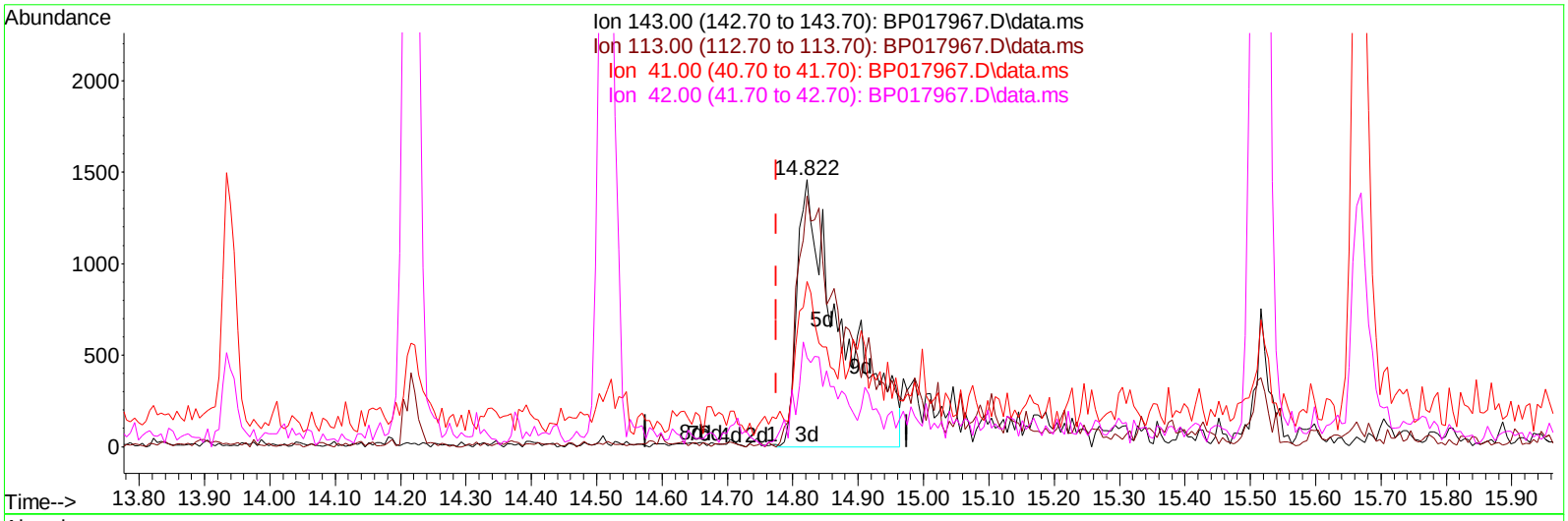

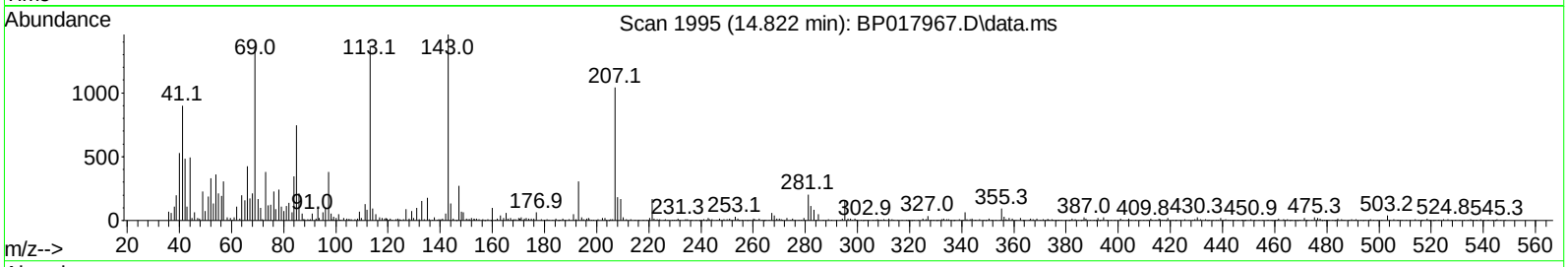

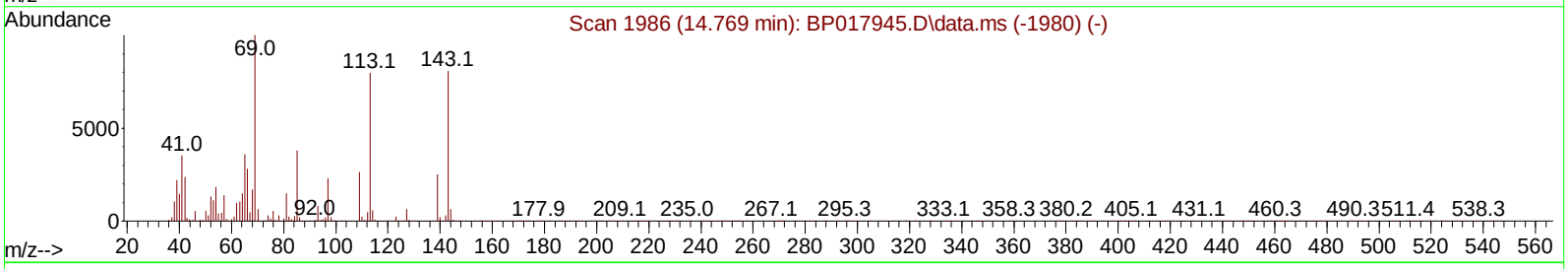

TIC: BP017967.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.047) 3.32 ng/ul m

response 6889

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	94.03
41.00	43.00	61.80#
42.00	28.20	33.33

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
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Quant Time: Nov 08 22:36:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.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	7.858	152	57406	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	243817	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.516	164	154104	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	341670	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	314676	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	344747	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	7164	4.776	ng/uL	0.00
4) Pyridine-d5	3.687	84	45122	10.840	ng/ul	0.01
7) Phenol-d5	7.028	99	26965	5.009	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	102490	31.561	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	94221	22.778	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	59052	13.272	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	64299	33.000	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	67217	29.952	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	111080	26.999	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	150218	25.196	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	468380	34.662	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	484571	32.681	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	6889m	3.316	ng/ul	0.05
60) Fluorene-d10	15.516	176	381723	33.751	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	64904	25.976	ng/ul	0.00
73) Anthracene-d10	17.392	188	626532	35.268	ng/ul	0.00
81) Pyrene-d10	19.639	212	791419	39.619	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	741888	38.514	ng/ul	0.00

Target Compounds Qvalue

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

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