Quantitation Report (Qedit)

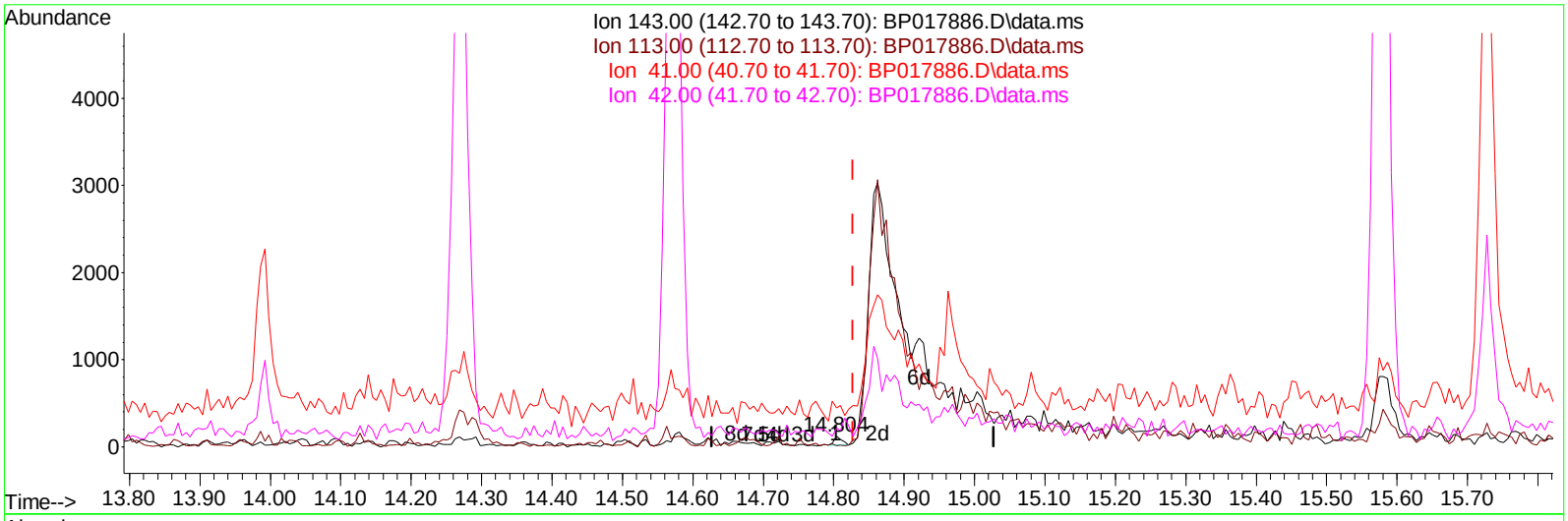
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017886.D
Acq On : 02 Nov 2023 14:58
Operator : MA/JU
Sample : 04996-16
Misc :
ALS Vial : 41 Sample Multiplier: 1

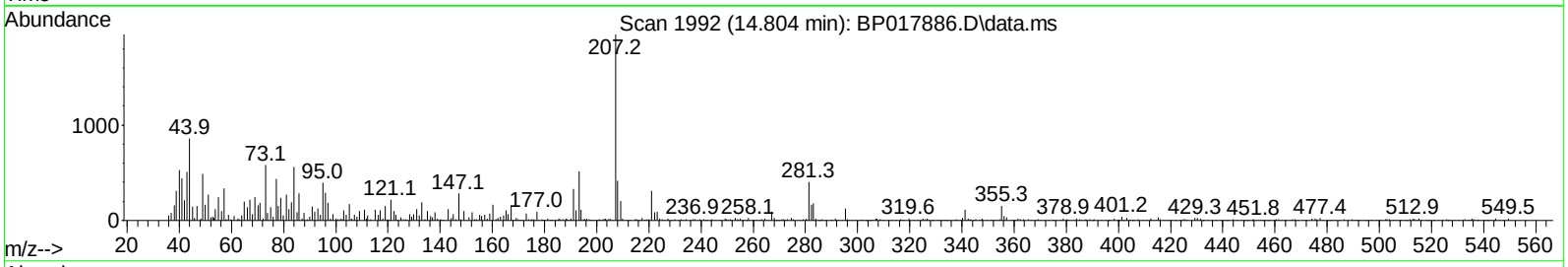
Quant Time: Nov 02 23:36:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 19 01:32:15 2023
Response via : Initial Calibration

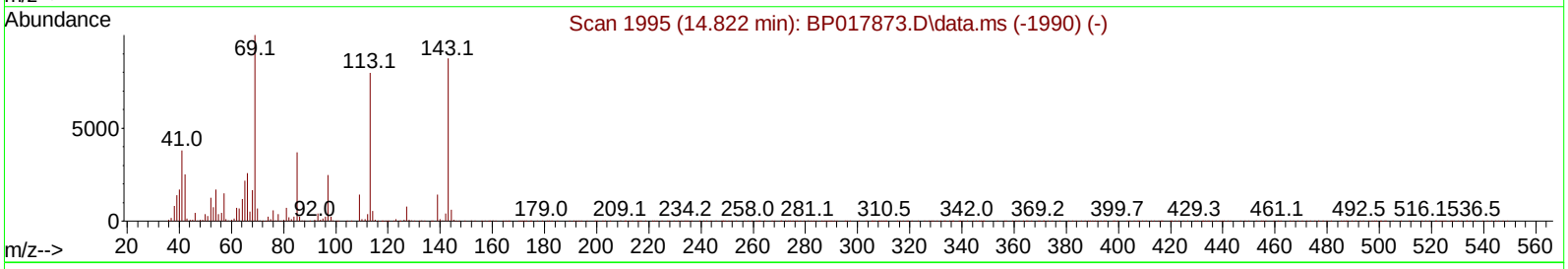
Instrument :
BNA_P
ClientSampleId :
A49C3

Manual IntegrationsAPPROVED

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







TIC: BP017886.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.804min (-0.024) 0.02 ng/ul

response 87

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	11.67#
41.00	39.20	371.67#
42.00	25.60	176.67#

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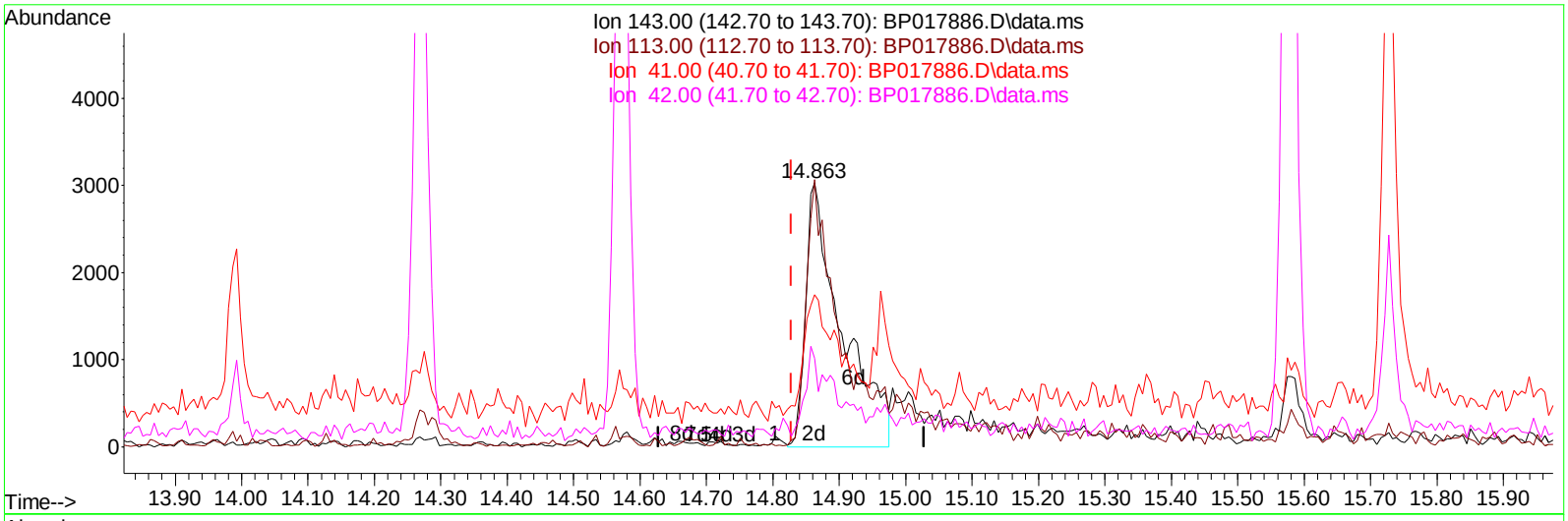

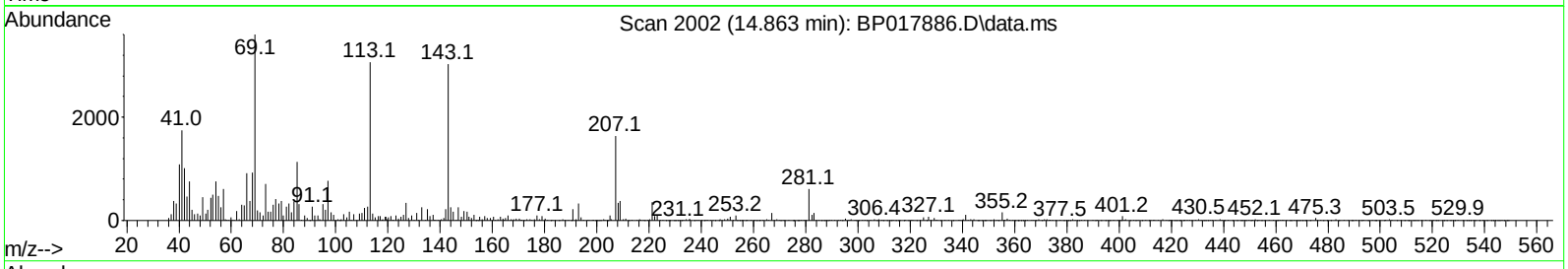

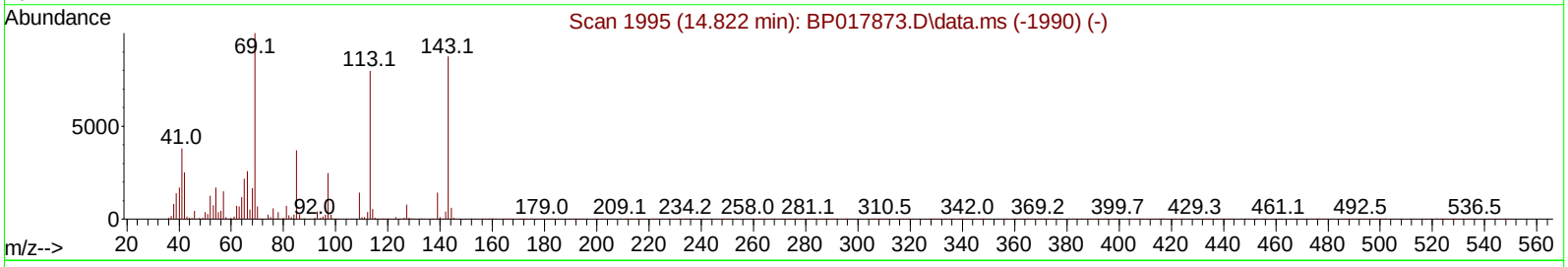

TIC: BP017886.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.035) 3.18 ng/ul m

response 11498

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	101.09#
41.00	39.20	57.57#
42.00	25.60	33.43#

Quantitation Report (QT Reviewed)

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Data File : BP017886.D
Acq On : 02 Nov 2023 14:58
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Quant Time: Nov 03 00:01:56 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	7.875	152	134475	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	513936	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	302308	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.368	188	622507	20.000	ng/ul	0.00
79) Chrysene-d12	21.509	240	522988	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264	534938	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	12320	3.408	ng/uL	0.01
4) Pyridine-d5	3.704	84	39377	4.161	ng/ul	0.01
7) Phenol-d5	7.045	99	50545	4.517	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	162207	23.471	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.398	132	165973	18.185	ng/ul	-0.01
15) 4-Methylphenol-d8	8.598	113	97309	11.073	ng/ul	-0.01
21) Nitrobenzene-d5	9.081	128	105496	25.332	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.798	143	115612	25.126	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	183491	20.329	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.881	131	217704	18.430	ng/ul	-0.01
46) Dimethylphthalate-d6	13.992	166	681627	26.859	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	712183	25.164	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.863	143	11498m	3.180	ng/ul	0.03
60) Fluorene-d10	15.580	176	554032	25.255	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.727	200	93517	22.807	ng/ul	0.00
73) Anthracene-d10	17.468	188	929925	29.232	ng/ul	0.00
81) Pyrene-d10	19.727	212	1096940	33.455	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	921864	30.941	ng/ul	0.00
Target Compounds						Qvalue

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

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