Quantitation Report (Qedit)

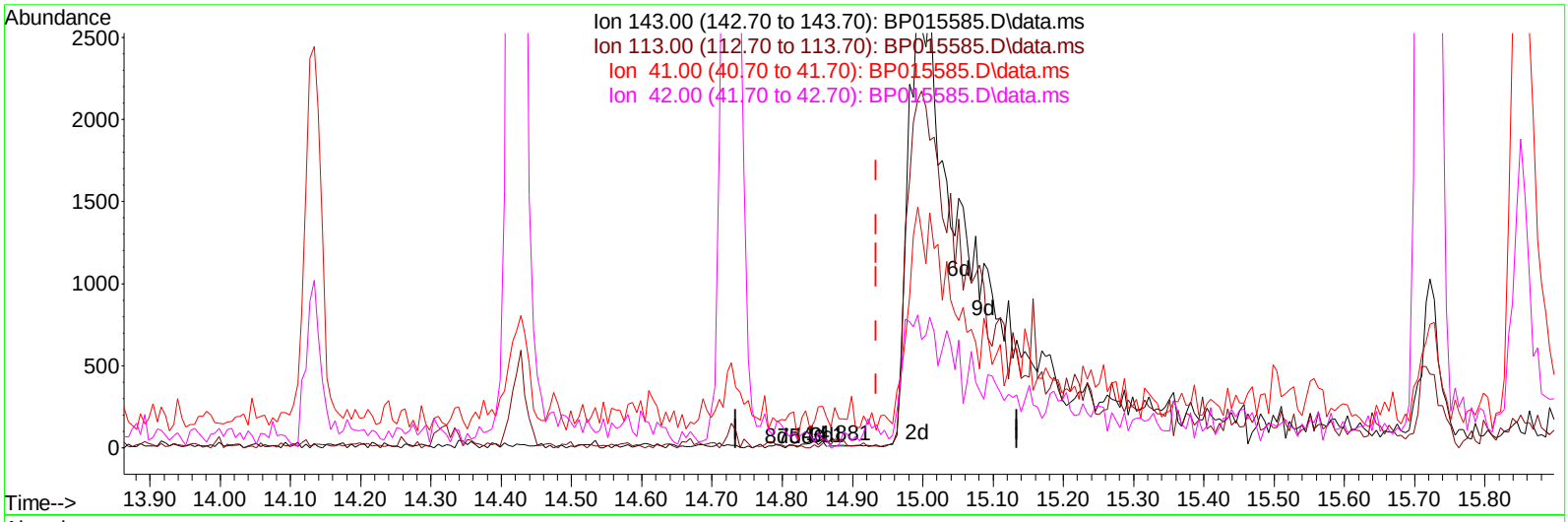
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015585.D
Acq On : 16 Jun 2023 17:29
Operator : MA/JU
Sample : 03220-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

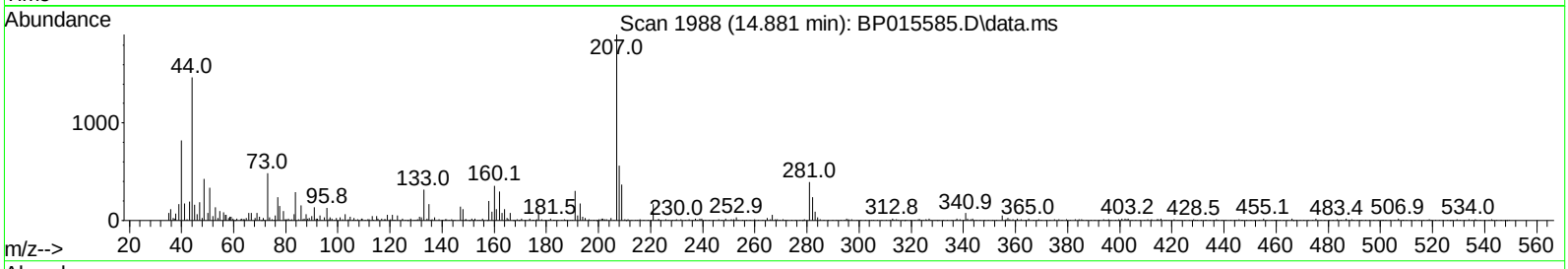
Quant Time: Jun 16 22:07:36 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 15 05:36:46 2023
Response via : Initial Calibration

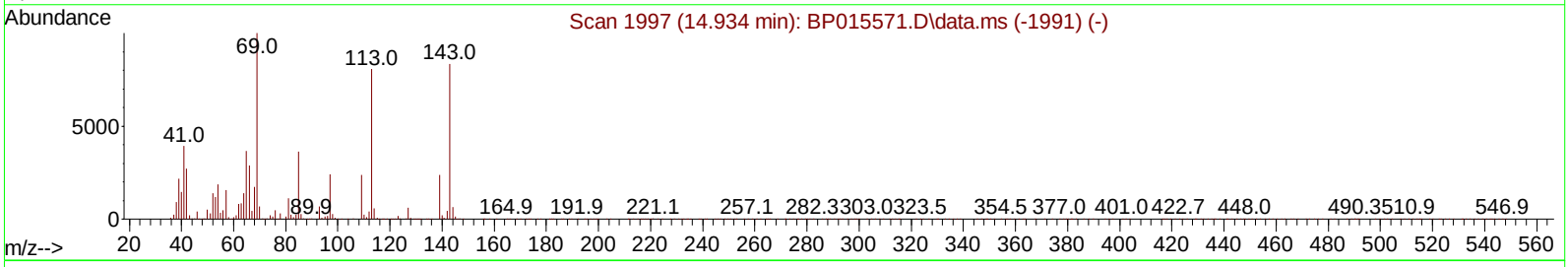
Instrument :
BNA_P
ClientSampleId :
C0NQ9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/19/2023
Supervised By :mohammad ahmed 06/20/2023







TIC: BP015585.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (-0.053) 0.00 ng/ul

response 14

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	346.67#
41.00	42.20	1160.00#
42.00	29.40	106.67#

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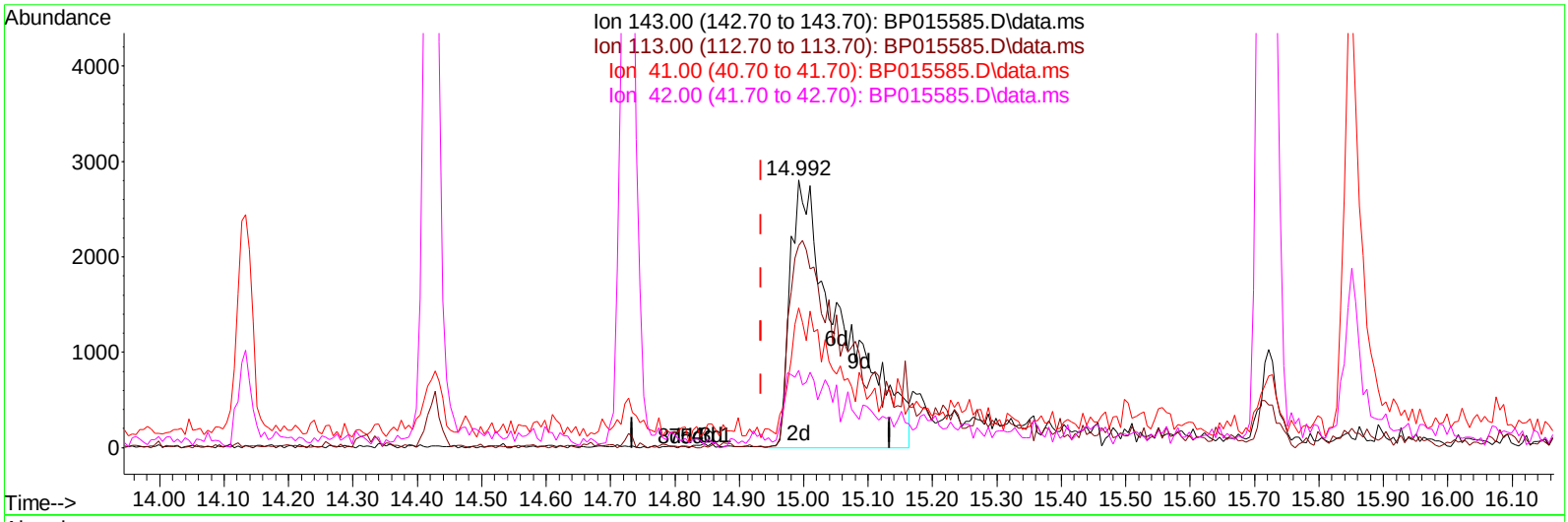

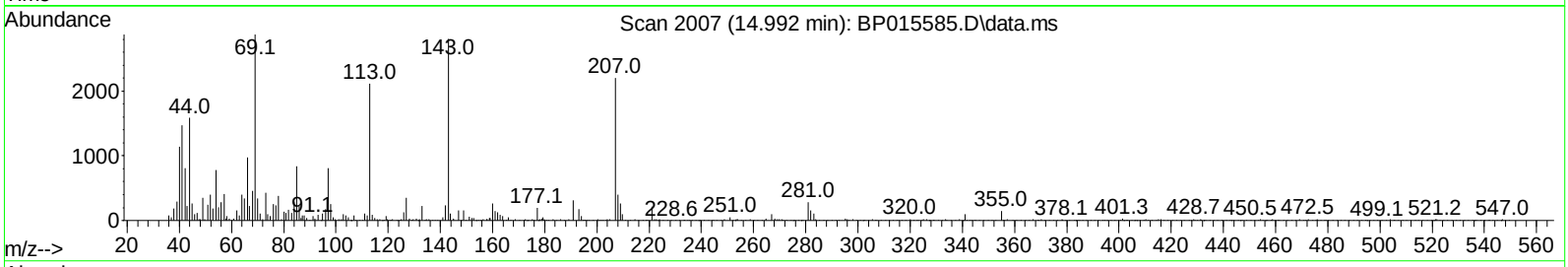

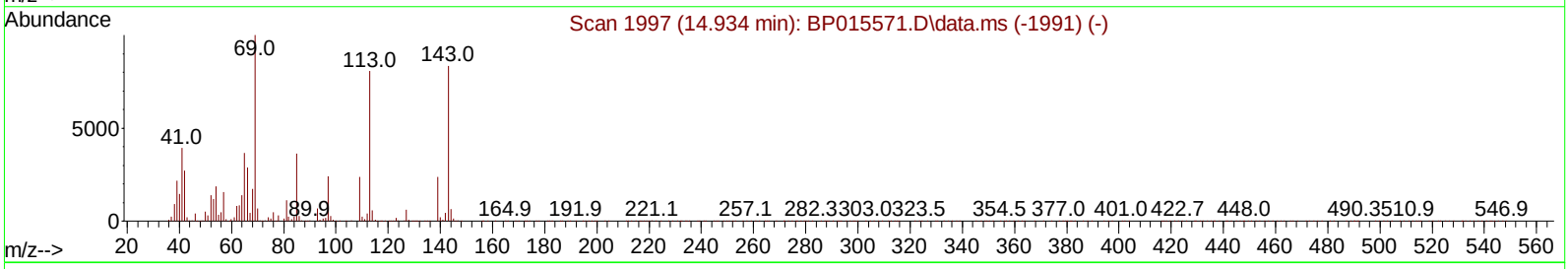

TIC: BP015585.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.992min (+ 0.059) 3.57 ng/ul m

response 15616

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	75.57
41.00	42.20	52.35#
42.00	29.40	28.92

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015585.D
Acq On : 16 Jun 2023 17:29
Operator : MA/JU
Sample : 03220-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Quant Time: Jun 16 22:59:28 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 15 05:36:46 2023
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	96091	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	427253	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	309331	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	752774	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	739844	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	826794	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	11926	4.596	ng/uL	0.00
4) Pyridine-d5	3.858	84	61382	9.107	ng/ul	0.01
7) Phenol-d5	7.246	99	51513	5.819	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	174111	32.933	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	157216	24.495	ng/ul	0.00
15) 4-Methylphenol-d8	8.799	113	108406	14.921	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	110713	33.006	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	116938	30.527	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	181533	25.979	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	260986	26.020	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	956317	39.205	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	942148	35.323	ng/ul	0.00
54) 4-Nitrophenol-d4	14.992	143	15616m	3.575	ng/ul	0.06
60) Fluorene-d10	15.722	176	806304	39.199	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	91928	19.277	ng/ul	0.00
73) Anthracene-d10	17.586	188	1439603	42.032	ng/ul	0.00
81) Pyrene-d10	19.810	212	1819961	45.961	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.945	264	1795231	43.274	ng/ul	0.00

Target Compounds Qvalue

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

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