Quantitation Report (Qedit)

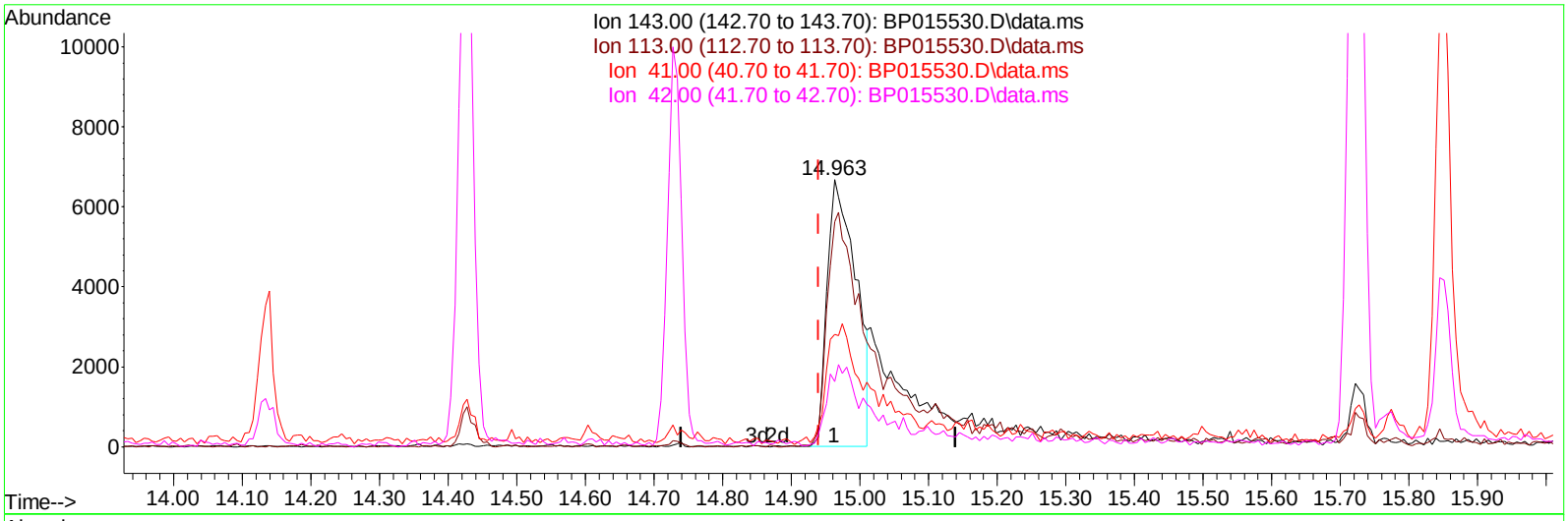
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015530.D
Acq On : 14 Jun 2023 22:15
Operator : MA/JU
Sample : 03104-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

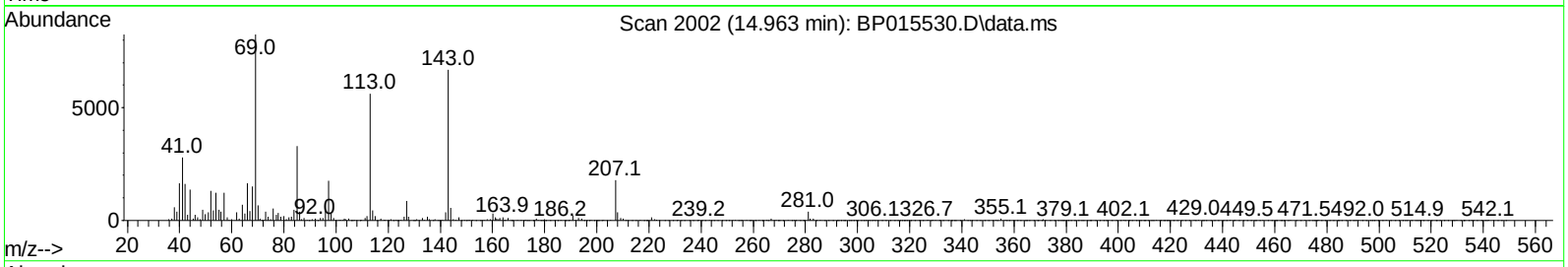
Quant Time: Jun 14 23:12:11 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

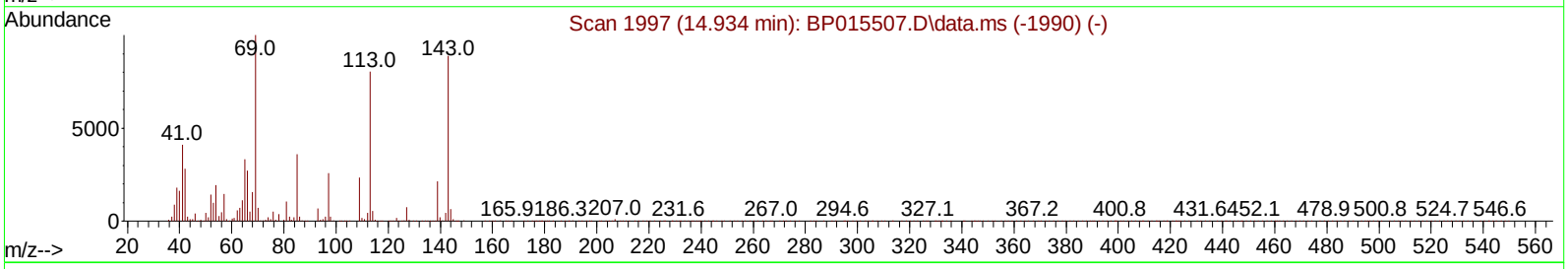
Instrument :
BNA_P
ClientSampleId :
C0NL5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023







TIC: BP015530.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 3.53 ng/ul

response 19374

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	84.27
41.00	42.20	42.02
42.00	29.40	24.34

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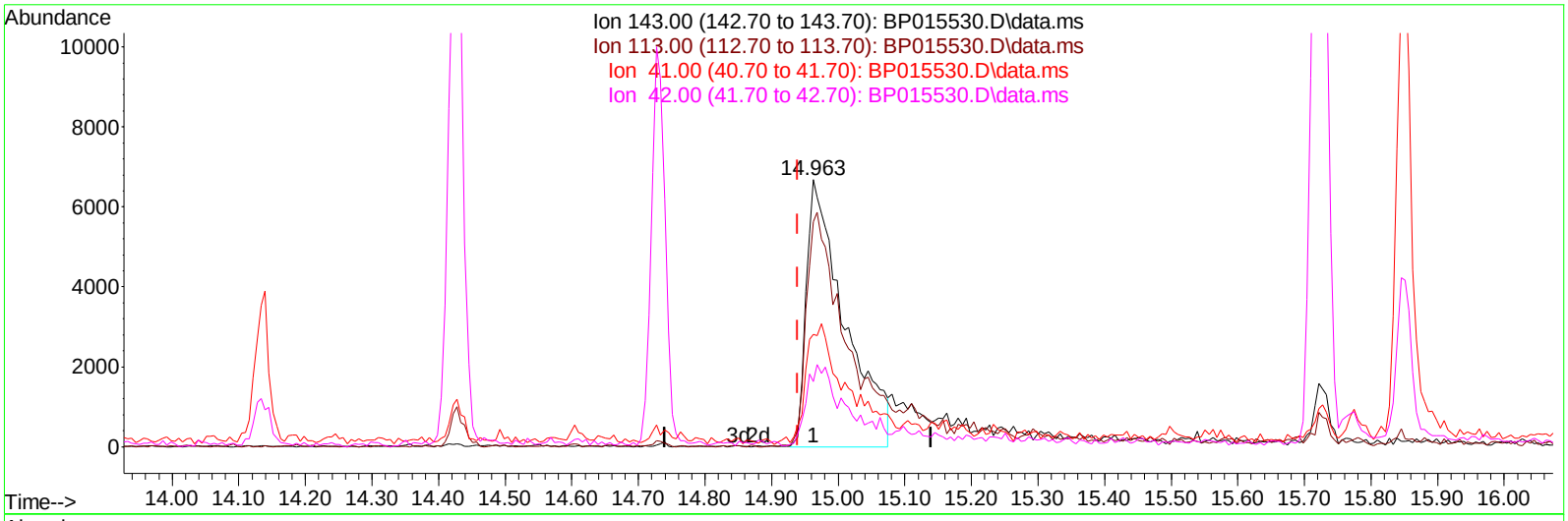

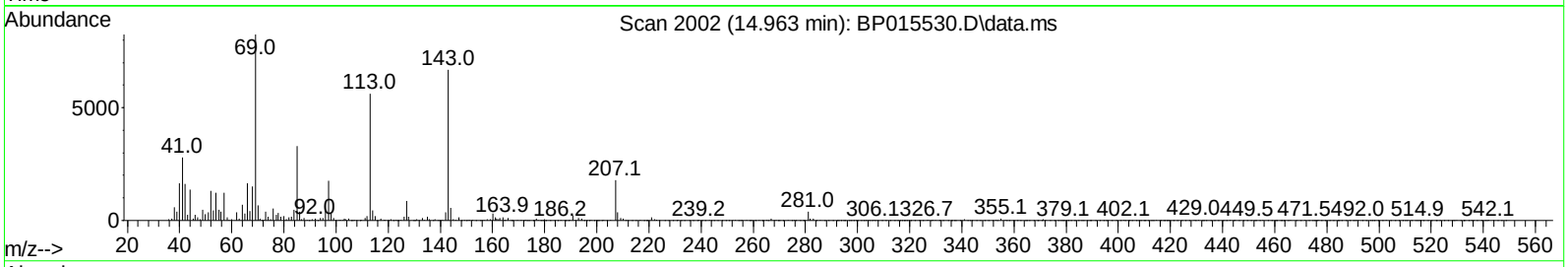

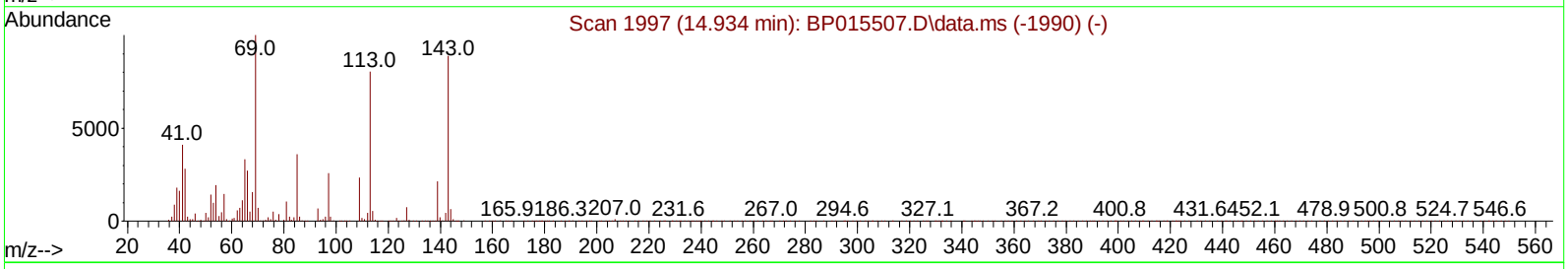

TIC: BP015530.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 4.87 ng/ul m

response 26680

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	84.27
41.00	42.20	42.02
42.00	29.40	24.34

Quantitation Report (QT Reviewed)

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Quant Time: Jun 15 00:03:01 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	120126	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	560893	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	388348	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	913508	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	948429	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1008051	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	20417	6.294	ng/uL	0.00
4) Pyridine-d5	3.857	84	88627	10.518	ng/ul	0.01
7) Phenol-d5	7.246	99	99792	9.017	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	279914	42.352	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	269976	33.647	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	197013	21.691	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	180866	41.073	ng/ul	0.00
24) 2-Nitrophenol-d4	9.992	143	206936	41.150	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	332474	36.243	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	335636	25.490	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	1318175	43.044	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1421125	42.440	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143	26680m	4.865	ng/ul	0.02
60) Fluorene-d10	15.722	176	1118448	43.310	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	216404	37.394	ng/ul	0.00
73) Anthracene-d10	17.586	188	1911053	45.979	ng/ul	0.00
81) Pyrene-d10	19.810	212	2376329	46.814	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2353299	46.526	ng/ul	0.00
Target Compounds						Qvalue

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

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