Quantitation Report (Qedit)

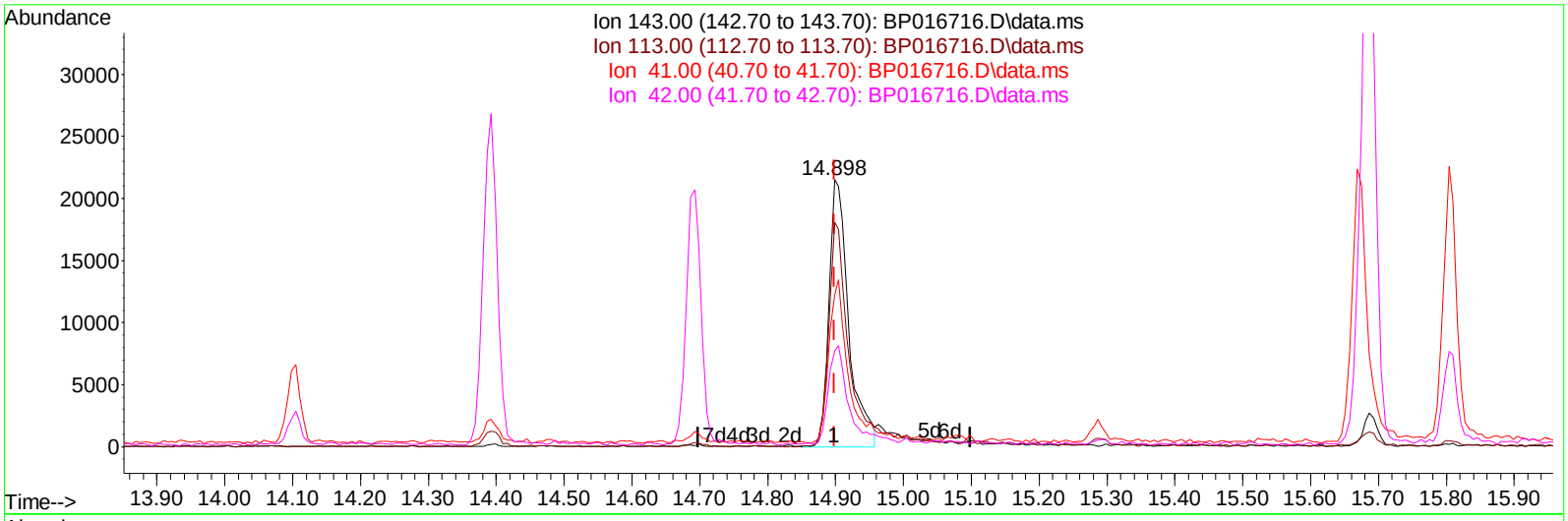
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082323\
Data File : BP016716.D
Acq On : 23 Aug 2023 15:50
Operator : MA/JU
Sample : 04105-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

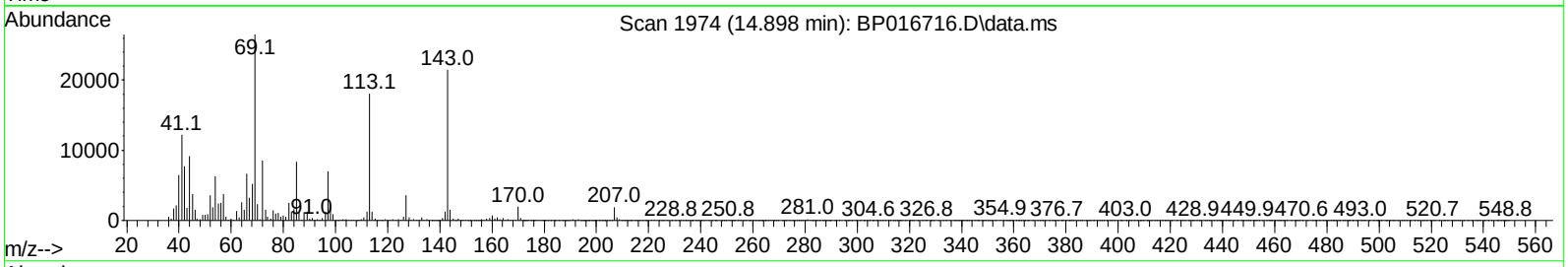
Quant Time: Aug 23 19:01:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 23 18:42:31 2023
Response via : Initial Calibration

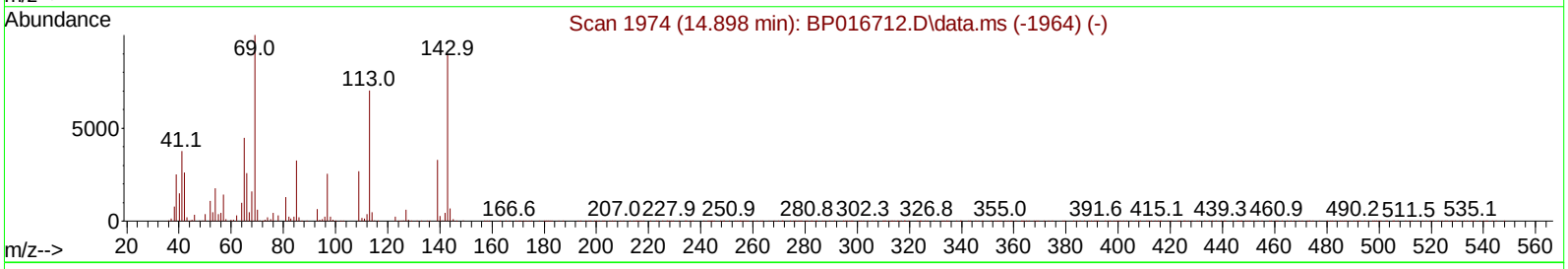
Instrument :
BNA_P
ClientSampleId :
BHAB8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/24/2023
Supervised By :mohammad ahmed 08/24/2023







TIC: BP016716.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.000) 3.56 ng/ul

response 43432

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	84.00
41.00	46.00	56.97#
42.00	30.30	35.90

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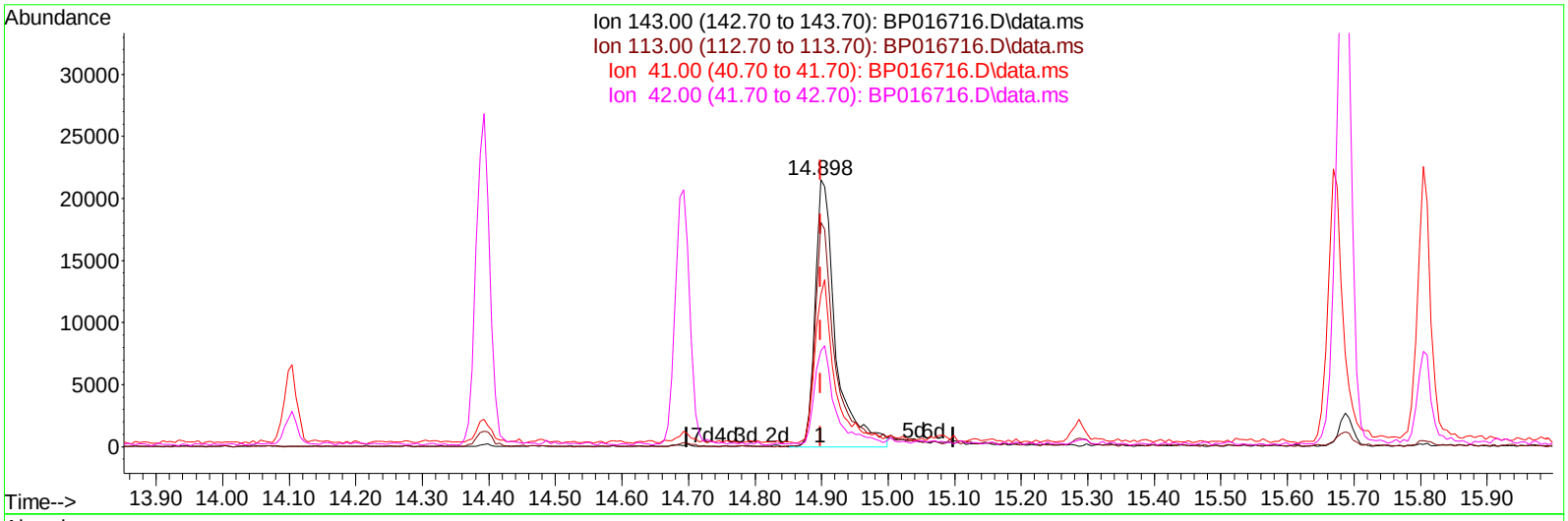

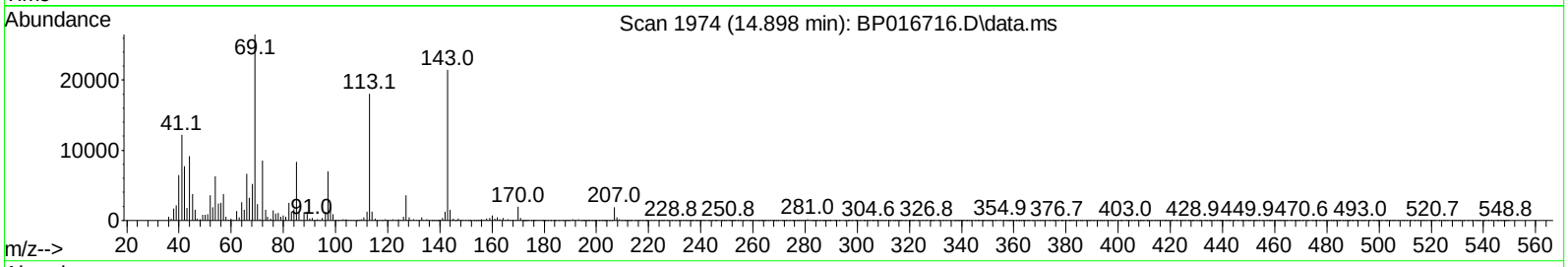

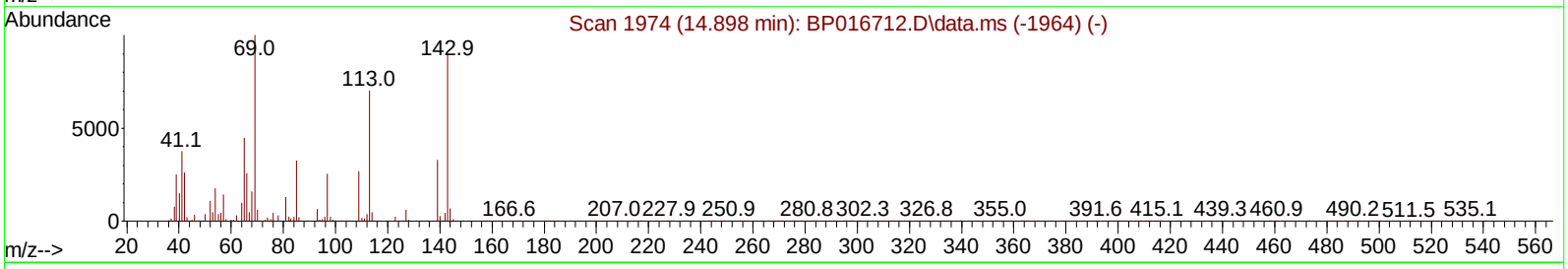

TIC: BP016716.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.000) 3.80 ng/ul m

response 46349

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	84.00
41.00	46.00	56.97#
42.00	30.30	35.90

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082323\
Data File : BP016716.D
Acq On : 23 Aug 2023 15:50
Operator : MA/JU
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Misc :
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Quant Time: Aug 23 19:07:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.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	8.040	152	288580	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	1338835	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.693	164	827553	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	1827516	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	1383510	20.000	ng/ul	0.00
88) Perylene-d12	24.133	264	1275224	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	46527	6.126	ng/uL	0.00
4) Pyridine-d5	3.828	84	218394	9.452	ng/ul	0.00
7) Phenol-d5	7.181	99	189854	7.136	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	664152	37.901	ng/ul	0.00
11) 2-Chlorophenol-d4	7.558	132	536344	28.068	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	362573	17.139	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	380851	40.428	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	369877	41.339	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	605347	34.379	ng/ul	0.00
31) 4-Chloroaniline-d4	11.028	131	952910	31.541	ng/ul	0.00
46) Dimethylphthalate-d6	14.104	166	2351042	39.278	ng/ul	0.00
49) Acenaphthylene-d8	14.393	160	2646303	39.170	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	46349m	3.801	ng/ul	0.00
60) Fluorene-d10	15.687	176	2048060	40.136	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	265376	30.524	ng/ul	0.00
73) Anthracene-d10	17.563	188	3240267	40.990	ng/ul	0.00
81) Pyrene-d10	19.792	212	3521557	48.578	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	2576856	42.165	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.423	88	29003	3.473	ng/uL	97

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

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