Quantitation Report (Qedit)

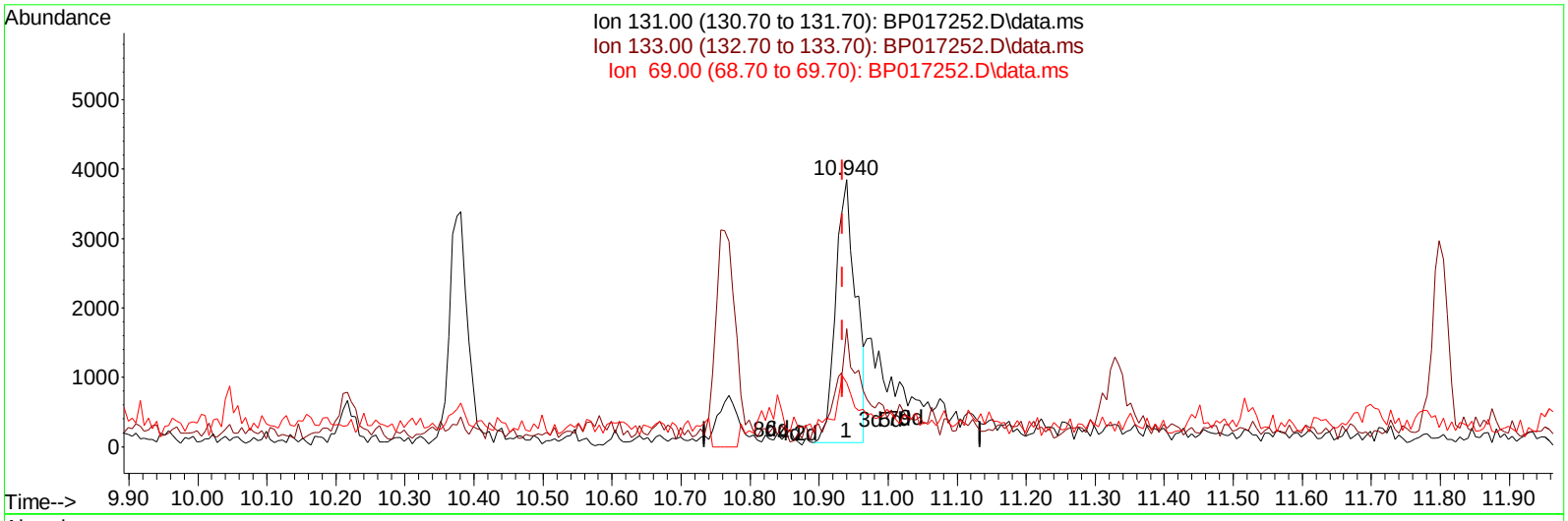
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017252.D
Acq On : 20 Sep 2023 22:34
Operator : MA/JU
Sample : 04332-16
Misc :
ALS Vial : 21 Sample Multiplier: 1

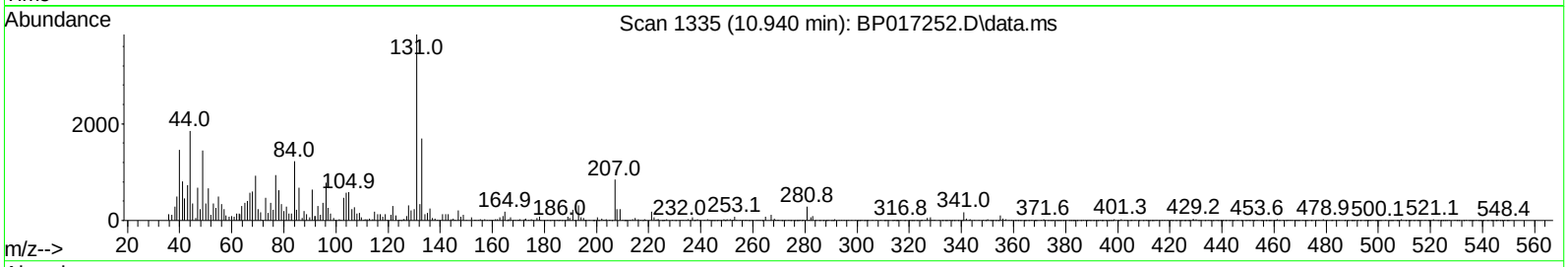
Quant Time: Sep 20 23:40:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 20 14:34:55 2023
Response via : Initial Calibration

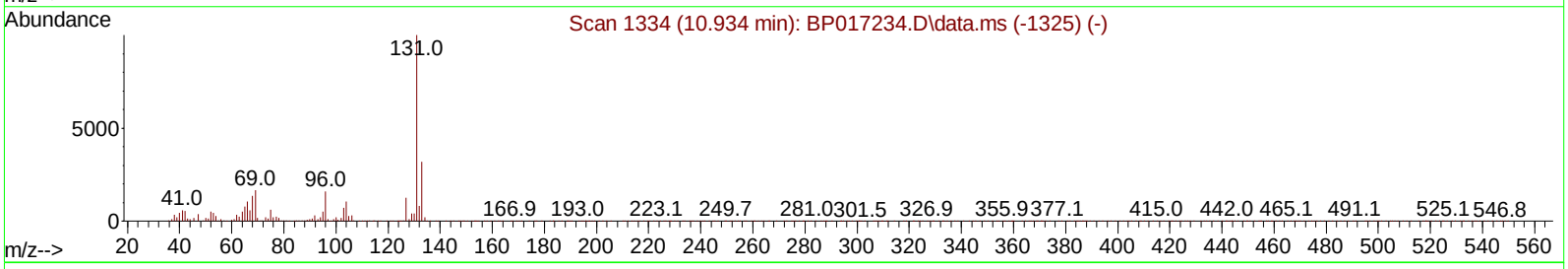
Instrument :
BNA_P
ClientSampleId :
C0AE3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/21/2023
Supervised By :mohammad ahmed 09/21/2023







TIC: BP017252.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.940min (+ 0.006) 0.53 ng/ul

response 7663

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	44.19#
69.00	16.70	24.02#
0.00	0.00	0.00

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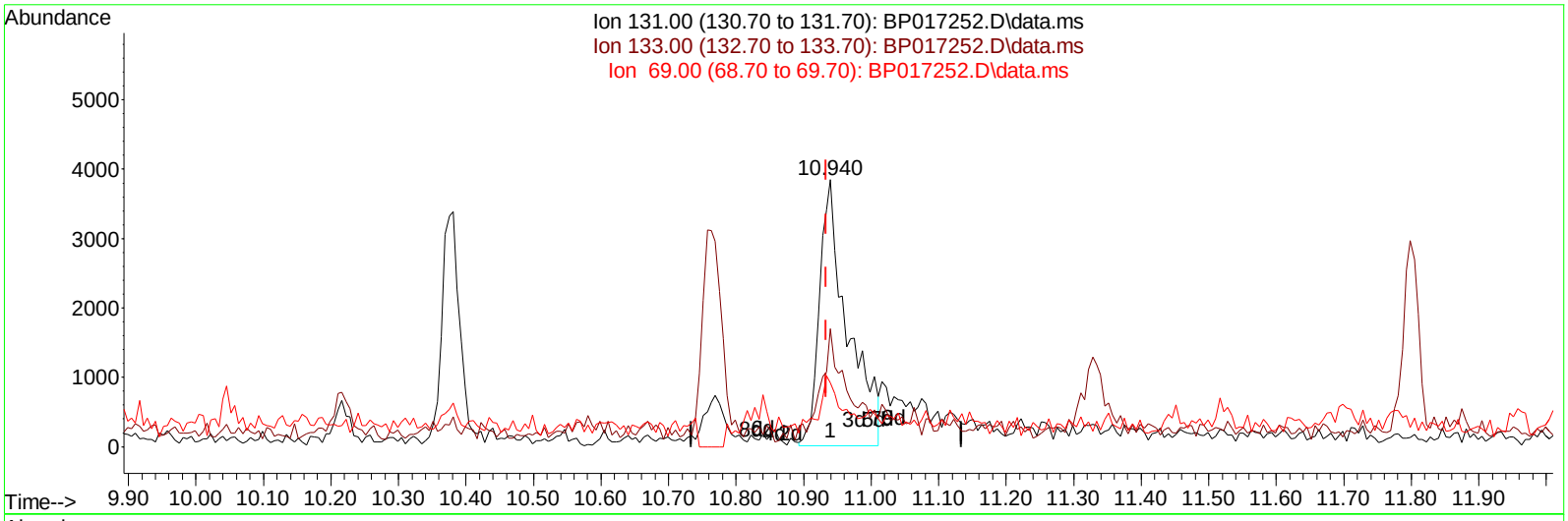

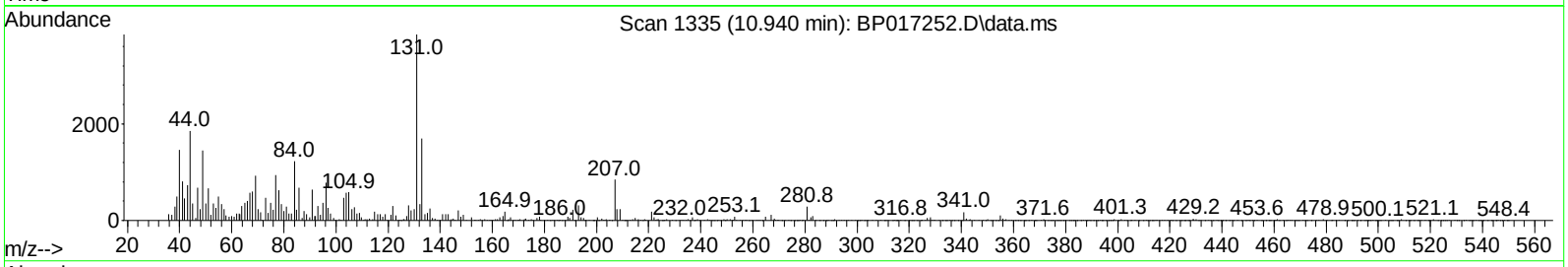

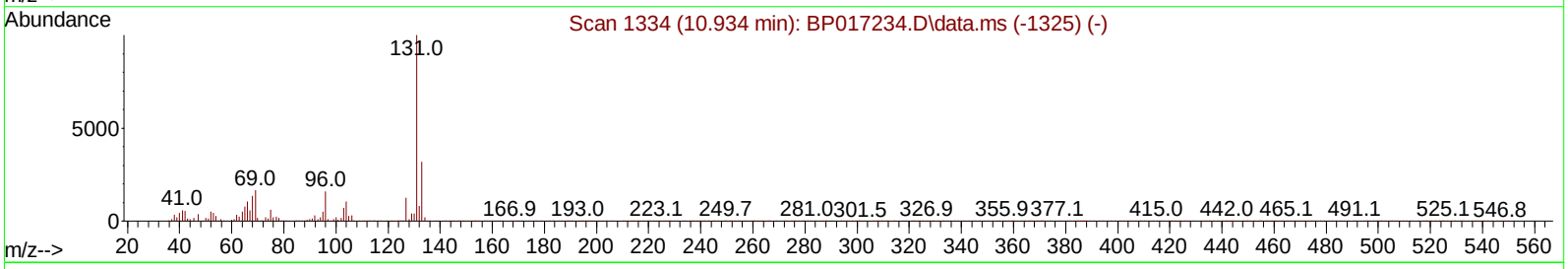

TIC: BP017252.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.940min (+ 0.006) 0.76 ng/ul m

response 11037

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	44.19#
69.00	16.70	24.02#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	173472	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	678632	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	398598	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	887756	20.000	ng/ul	-0.01
79) Chrysene-d12	21.475	240	826822	20.000	ng/ul	-0.01
88) Perylene-d12	23.998	264	977722	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	13124	3.107	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.093	99	224396	15.750	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	143996	16.747	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	189969	16.812	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	167959	15.163	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	91606	18.672	ng/ul	0.00
24) 2-Nitrophenol-d4	9.852	143	101513	19.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	182999	18.183	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	11037m	0.757	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	559137	19.582	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	639162	19.462	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	66189	13.305	ng/ul	0.00
60) Fluorene-d10	15.598	176	491897	20.010	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	63323	12.656	ng/ul	0.00
73) Anthracene-d10	17.469	188	784170	19.917	ng/ul	-0.01
81) Pyrene-d10	19.710	212	918692	20.547	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.827	264	906452	19.256	ng/ul	-0.02
Target Compounds						Qvalue
6) Benzaldehyde	7.099	77	92649	12.743	ng/ul	94
8) Phenol	7.122	94	43742	3.053	ng/ul	95
16) Acetophenone	8.787	105	250198	14.617	ng/ul	98

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

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