

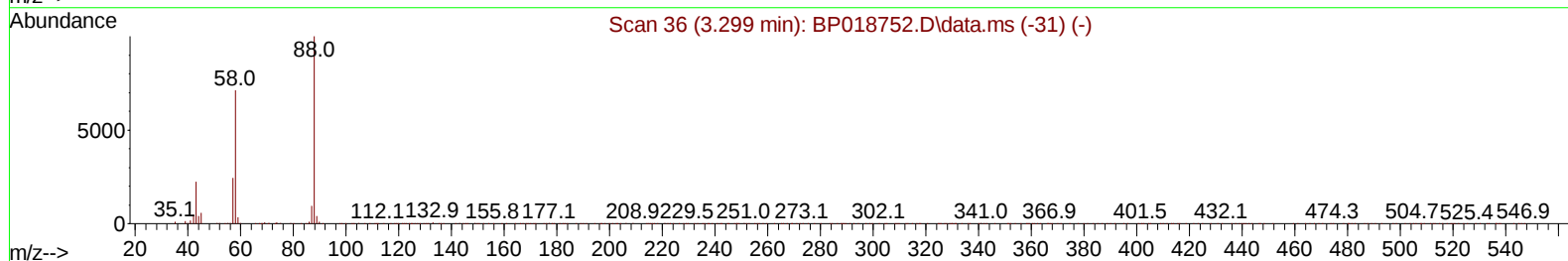
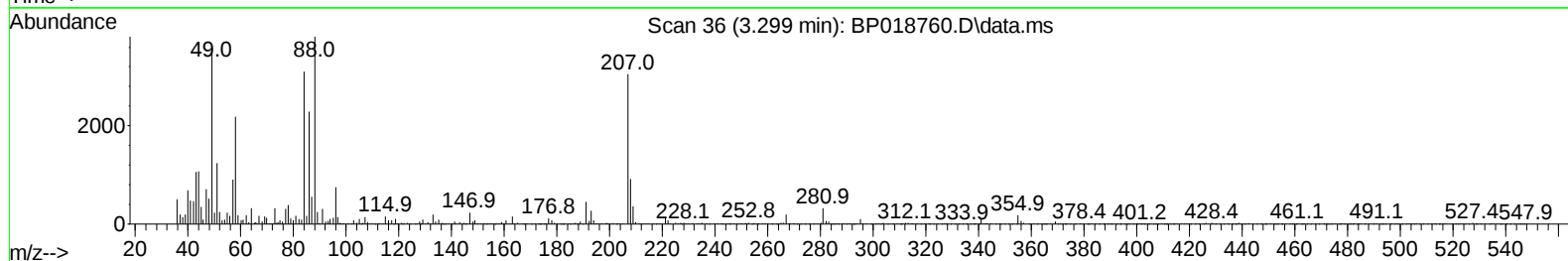
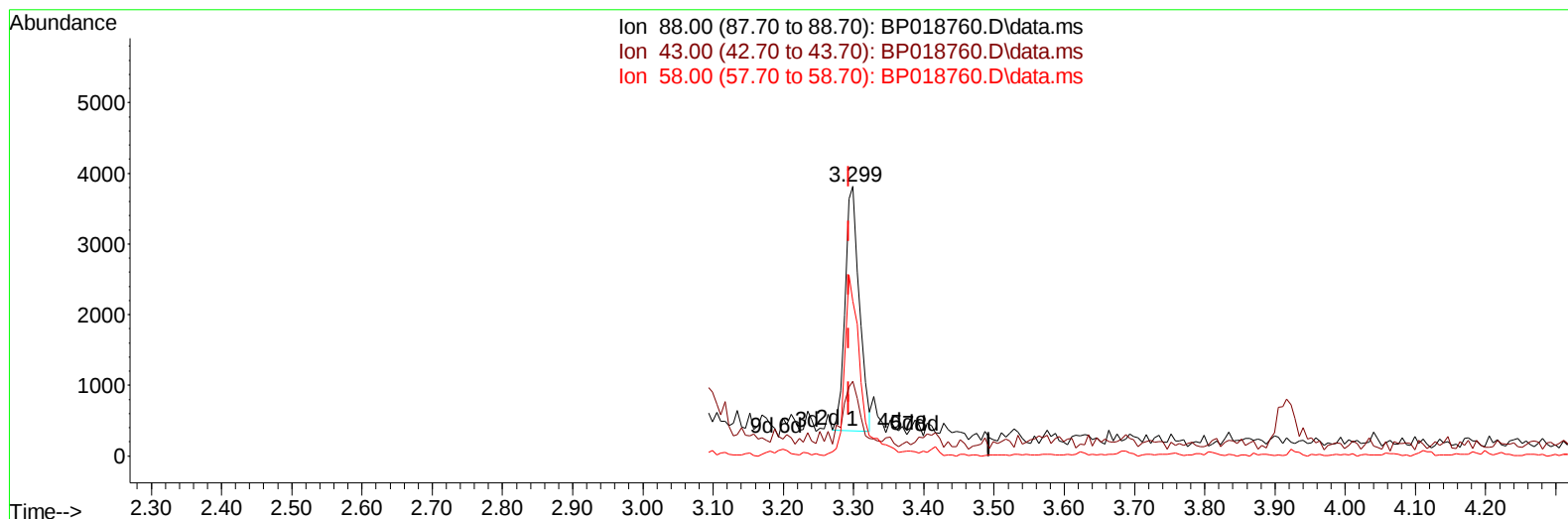
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018760.D
Acq On : 12 Dec 2023 10:39
Operator : MA/JU
Sample : 05675-10
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0AB1

Manual IntegrationsAPPROVED

Quant Time: Dec 12 21:36:24 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 11 22:02:13 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/13/2023
Supervised By :mohammad ahmed 12/14/2023



TIC: BP018760.D\data.ms

(2) 1,4-Dioxane

3.299min (+ 0.006) 1.60 ng/uL

response 4857

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.20	27.77
58.00	70.60	57.44
0.00	0.00	0.00

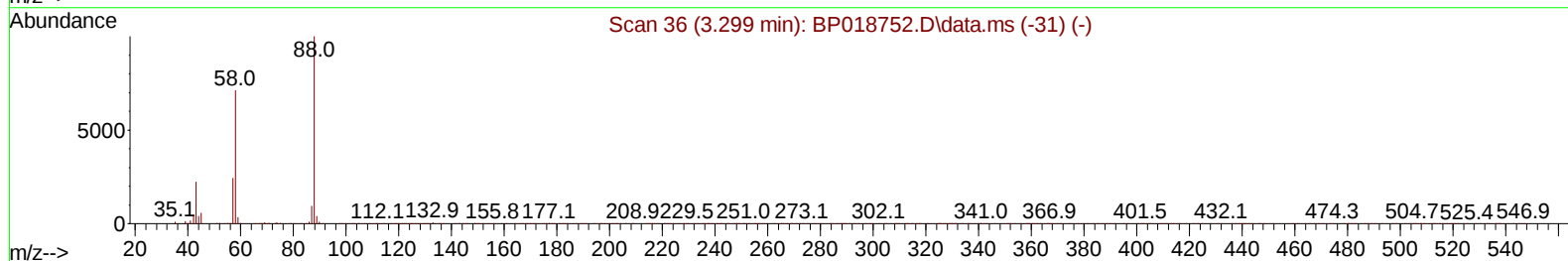
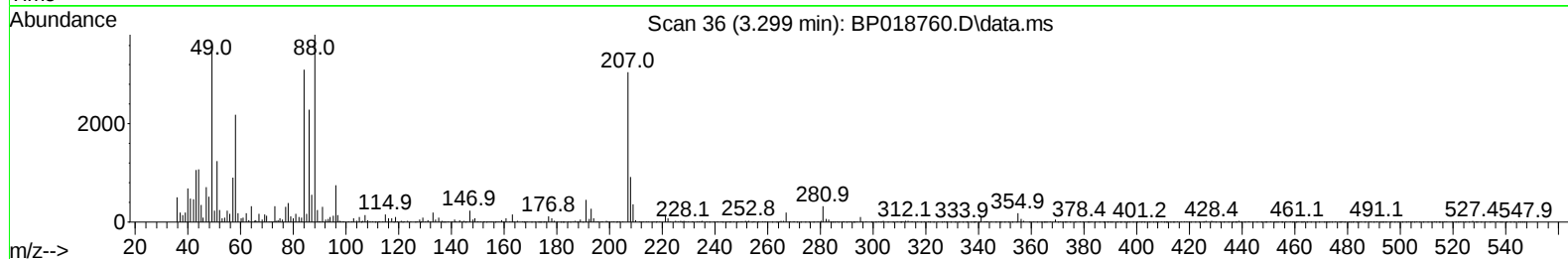
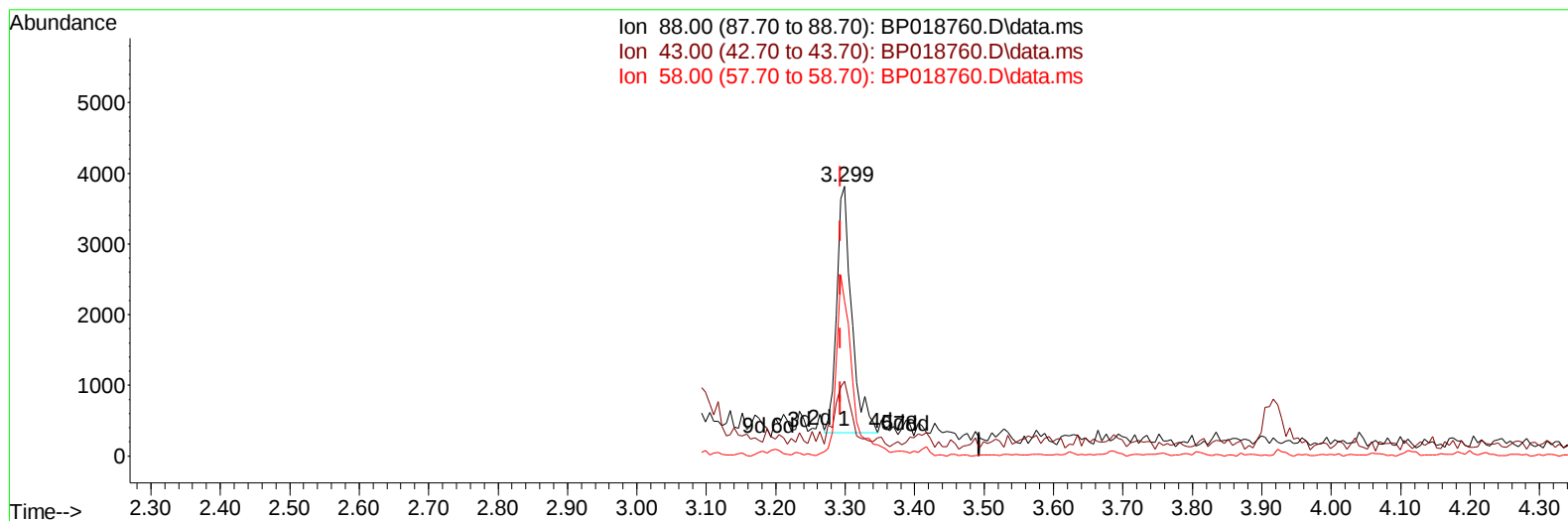
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TIC: BP018760.D\data.ms

(2) 1,4-Dioxane

3.299min (+ 0.006) 1.73 ng/uL m

response 5281

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.20	27.77
58.00	70.60	57.44
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/13/2023
 Supervised By :mohammad ahmed 12/14/2023

Quant Time: Dec 12 21:53:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 11 22:02:13 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	94623	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	338902	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	205223	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	475695	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.304	240	535043	20.000	ng/ul	# 0.00
88) Perylene-d12	23.663	264	637621	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	10207	3.676	ng/uL	0.00
4) Pyridine-d5	3.681	84	65468	8.802	ng/ul	0.00
7) Phenol-d5	7.011	99	38081	4.010	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	131747	23.409	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	129917	17.534	ng/ul	0.00
15) 4-Methylphenol-d8	8.552	113	70857	9.220	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	81267	26.942	ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	81479	26.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.258	165	151759	23.733	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	177472	21.237	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	553670	30.228	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	593104	29.407	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	10935	3.637	ng/ul	0.00
60) Fluorene-d10	15.457	176	489348	31.828	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	67437	23.643	ng/ul	0.00
73) Anthracene-d10	17.316	188	784263	31.360	ng/ul	0.00
81) Pyrene-d10	19.551	212	1013195	24.327	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	1173844	31.574	ng/ul	-0.02
Target Compounds						
2) 1,4-Dioxane	3.299	88	5281m	1.734	ng/uL	Qvalue

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

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