

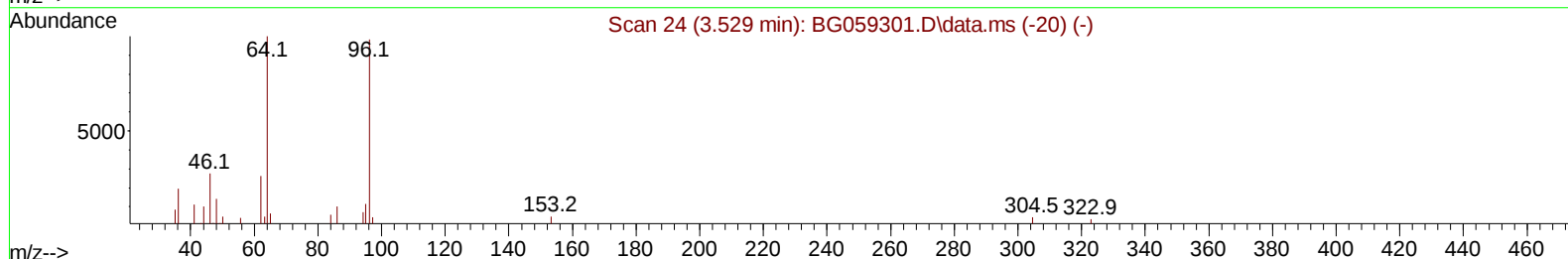
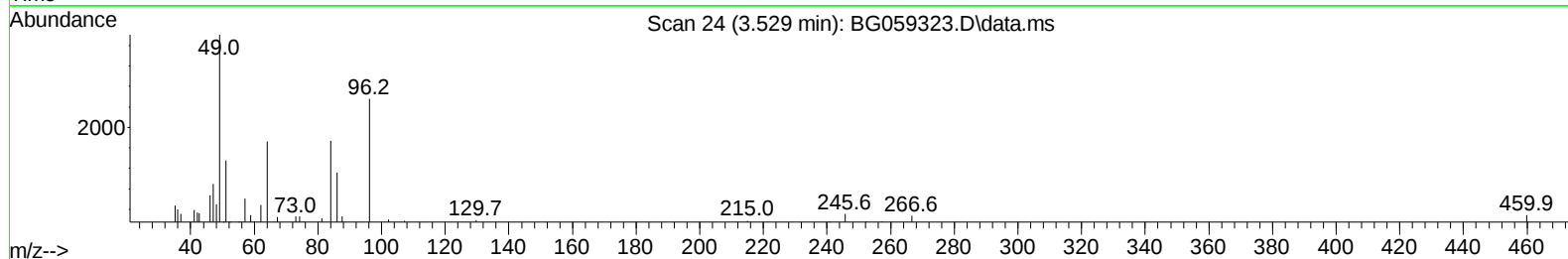
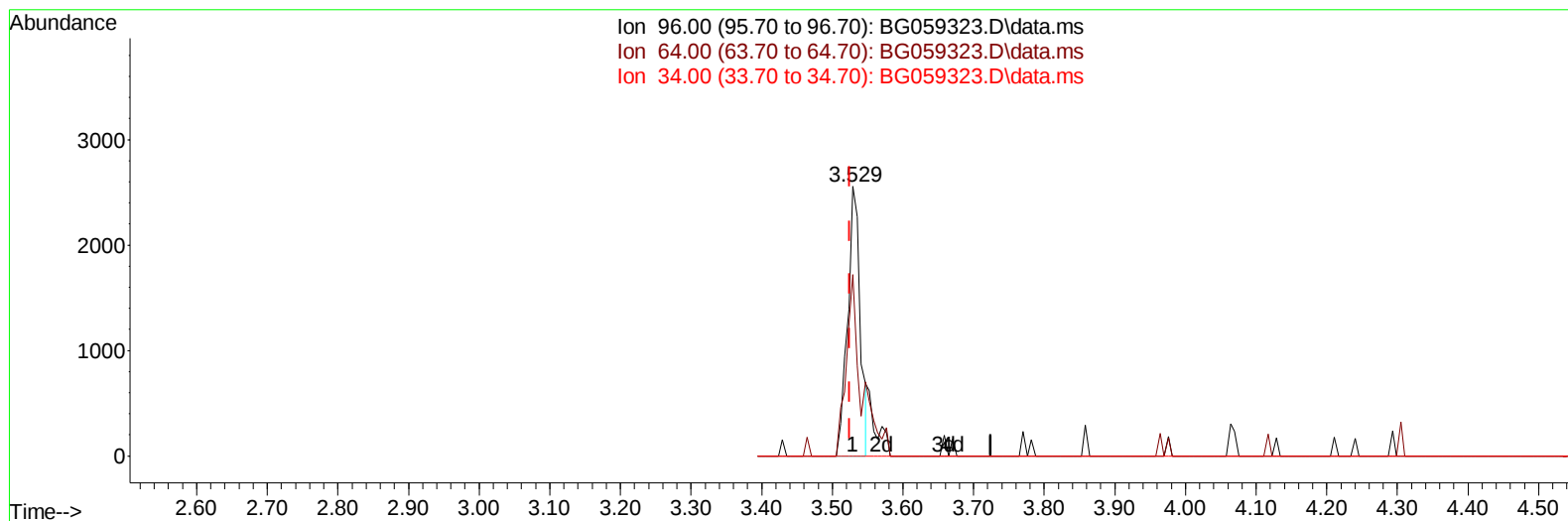
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
Operator : MA/JU
Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZY2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/07/2023
Supervised By :mohammad ahmed 10/07/2023

Quant Time: Oct 06 22:28:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 28 04:33:25 2023
Response via : Initial Calibration



TIC: BG059323.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.529min (+ 0.004) 2.59 ng/uL

response 3182

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	79.20	67.29
34.00	0.00	0.00
0.00	0.00	0.00

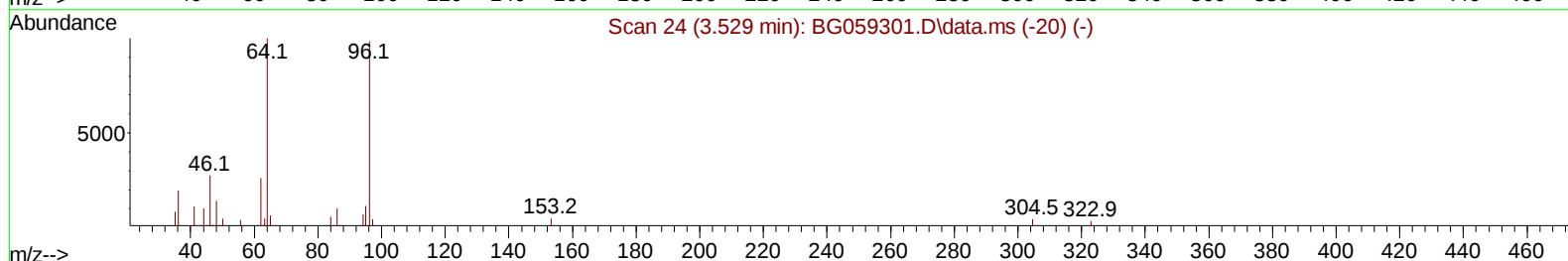
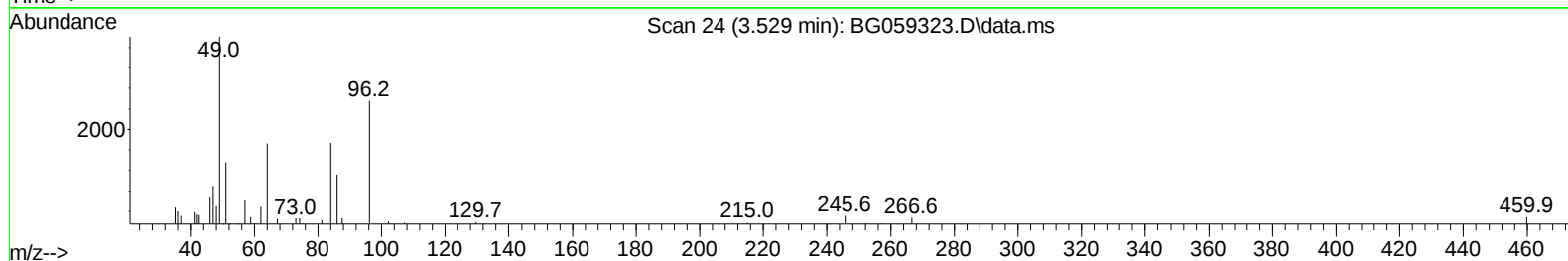
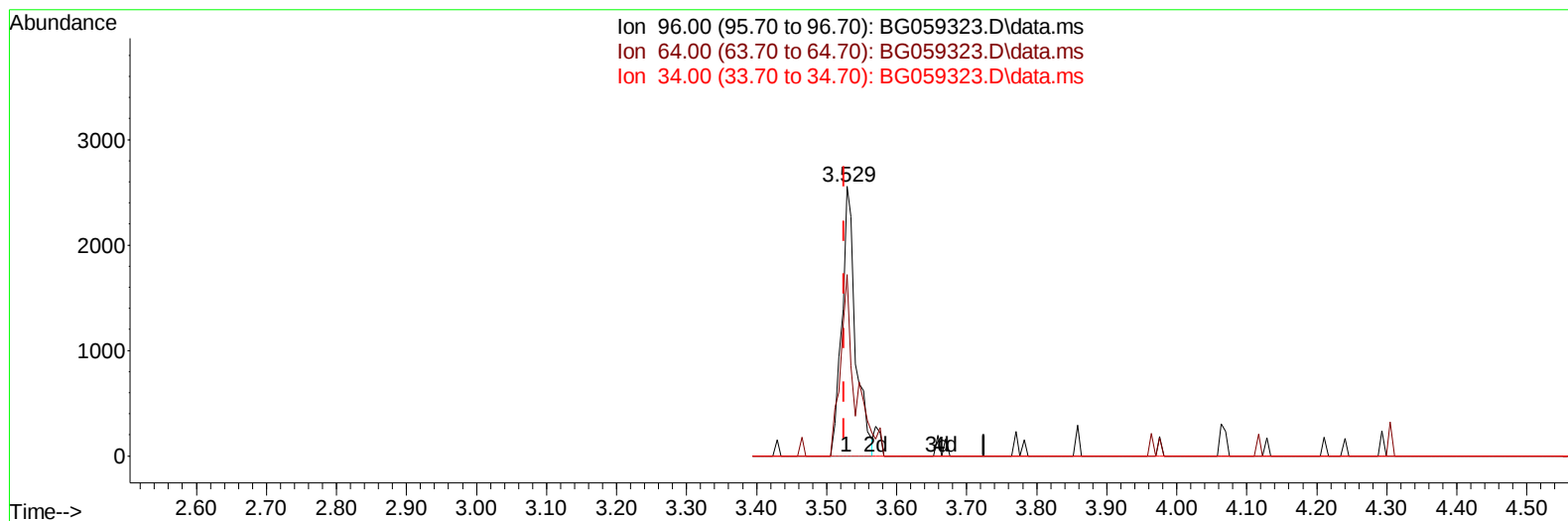
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TIC: BG059323.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.529min (+ 0.004) 2.88 ng/uL m

response 3538

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	79.20	67.29
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	46445	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	219306	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.863	164	138743	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.619	188	311737	20.000	ng/ul	0.00
79) Chrysene-d12	21.914	240	290750	20.000	ng/ul	0.00
88) Perylene-d12	25.404	264	338244	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	3538m	2.875	ng/uL	0.00
4) Pyridine-d5	3.964	84	41148	11.460	ng/ul	0.00
7) Phenol-d5	7.360	99	32762	7.047	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.554	67	100727	33.916	ng/ul	0.00
11) 2-Chlorophenol-d4	7.748	132	96601	30.911	ng/ul	0.00
15) 4-Methylphenol-d8	8.923	113	61327	17.065	ng/ul	0.00
21) Nitrobenzene-d5	9.428	128	65085	42.057	ng/ul	0.00
24) 2-Nitrophenol-d4	10.151	143	70497	46.403	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.680	165	115877	37.129	ng/ul	0.00
31) 4-Chloroaniline-d4	11.220	131	160225	31.732	ng/ul	0.00
46) Dimethylphthalate-d6	14.258	166	402778	40.153	ng/ul	0.00
49) Acenaphthylene-d8	14.569	160	495556	39.359	ng/ul	0.00
54) 4-Nitrophenol-d4	15.057	143	11979	6.839	ng/ul	0.00
60) Fluorene-d10	15.856	176	382339	40.047	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.968	200	52946	33.042	ng/ul	0.00
73) Anthracene-d10	17.713	188	595587	41.930	ng/ul	0.00
81) Pyrene-d10	19.986	212	695188	42.696	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.163	264	697689	42.923	ng/ul	0.00

Target Compounds Qvalue

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

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