

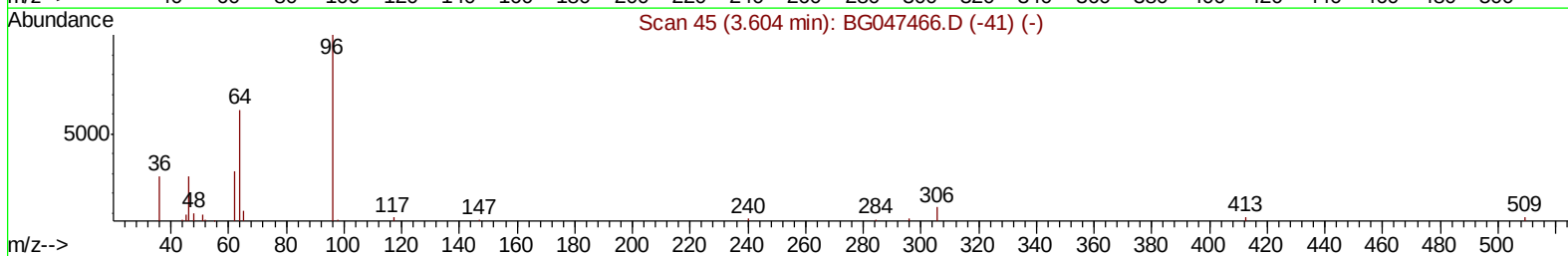
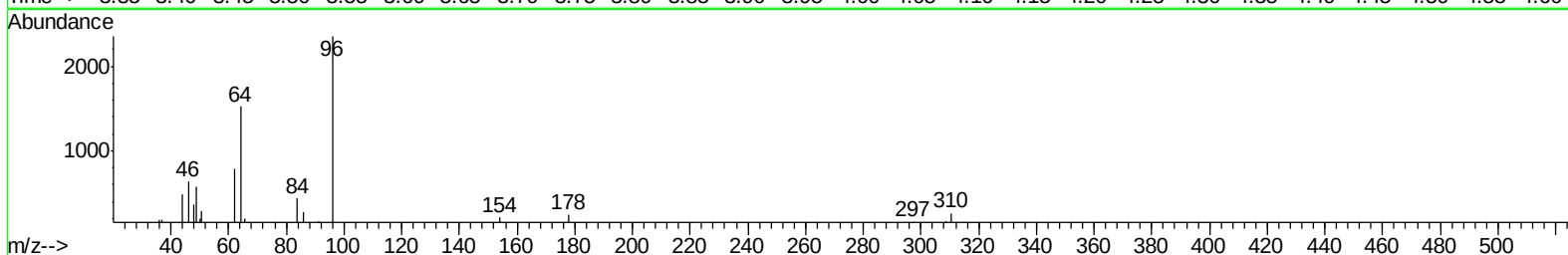
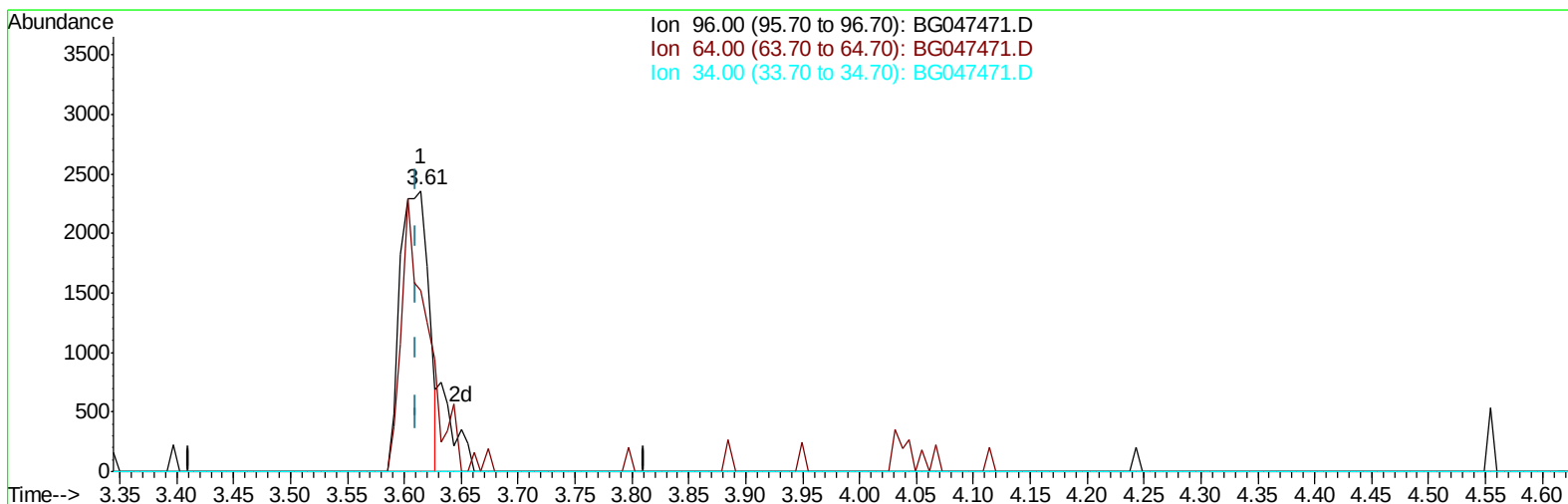
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG112520\
Data File : BG047471.D
Acq On : 25 Nov 2020 17:00
Operator : CG/JU
Sample : PB133233BL
Misc : MDL-WATER-BL03
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK233

Manual Integrations
APPROVED

mohammad
11/27/2020 11:45:47 AM

Quant Time: Nov 25 17:42:15 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG111820.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 25 12:54:28 2020
Response via : Initial Calibration



TIC: BG047471.D

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

3.614min (+0.004) 5.35ng/uL

response 4108

Ion	Exp%	Act%
96.00	100	100
64.00	61.30	64.70
34.00	0.00	0.00
0.00	0.00	0.00

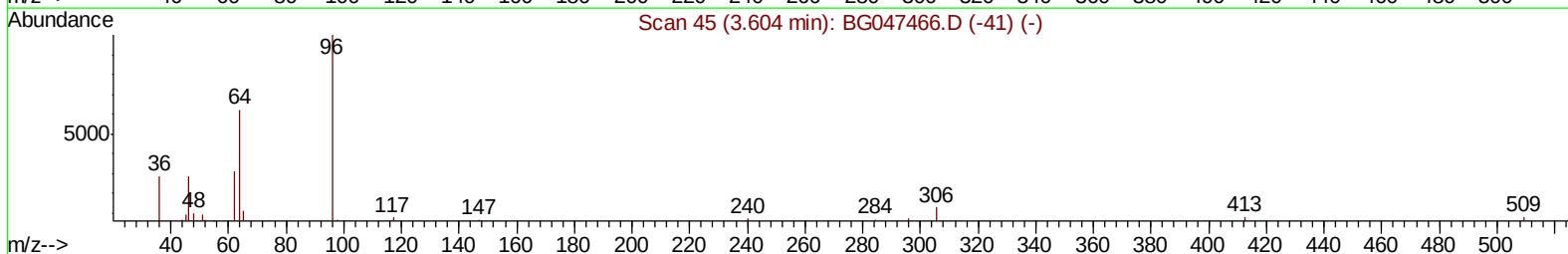
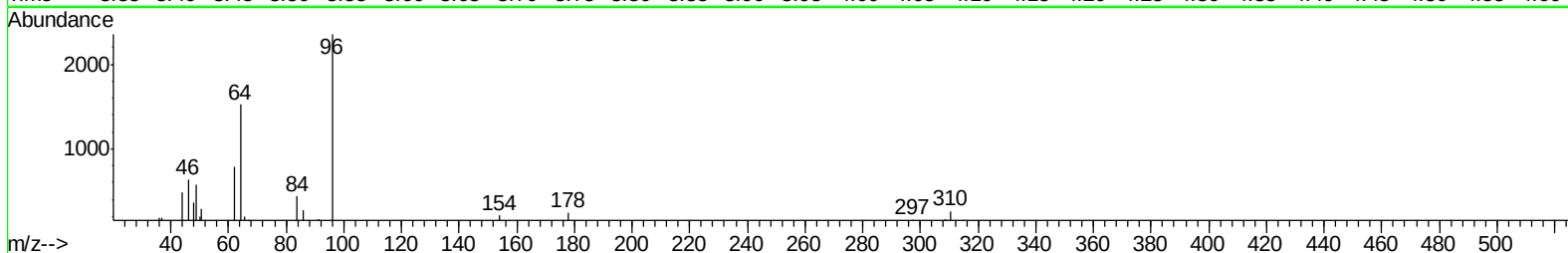
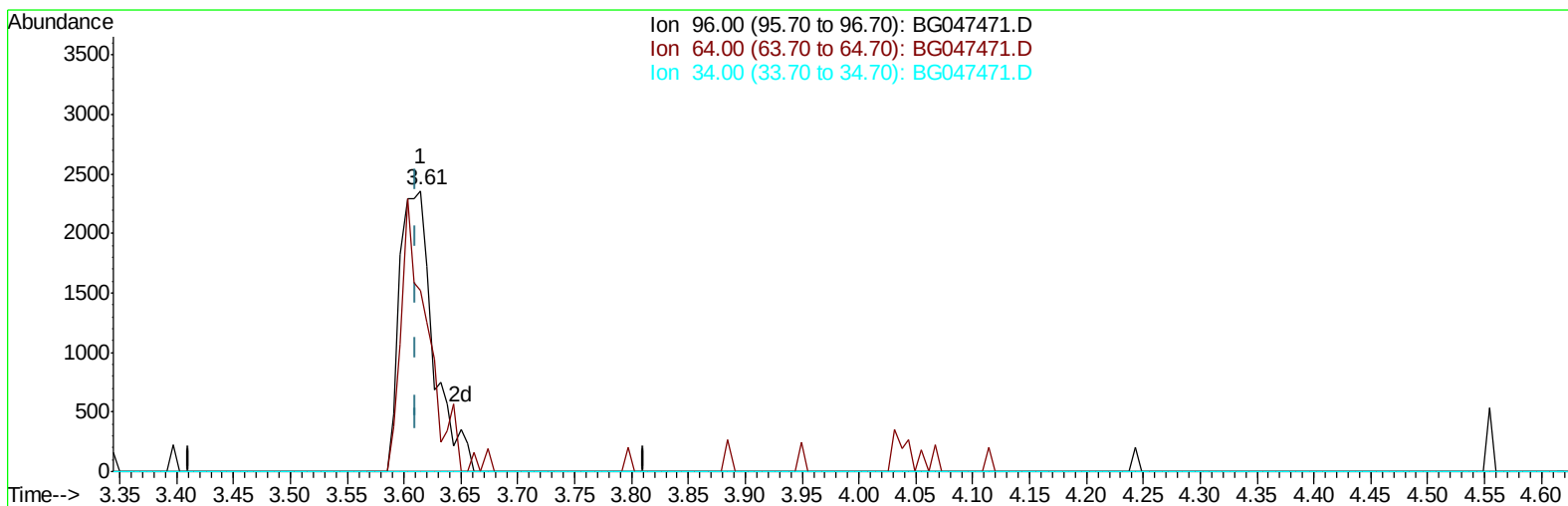
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TIC: BG047471.D

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

3.614min (+0.004) 6.33ng/uL m

response 4860

Ion	Exp%	Act%
96.00	100	100
64.00	61.30	64.70
34.00	0.00	0.00
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	30418	20.00	ng/ul	0.00
20) Naphthalene-d8	11.09	136	116405	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.88	164	90877	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	226283	20.00	ng/ul	0.00
79) Chrysene-d12	21.90	240	233649	20.00	ng/ul	0.00
88) Perylene-d12	25.29	264	237393	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.61	96	4860m	6.33	ng/uL	0.00
4) Pyridine-d5	4.04	84	76184	32.32	ng/ul	0.00
7) Phenol-d5	7.40	99	102011	35.00	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.58	67	59101	34.65	ng/ul	0.00
11) 2-Chlorophenol-d4	7.79	132	76439	38.43	ng/ul	0.00
15) 4-Methylphenol-d8	8.96	113	80388	35.51	ng/ul	0.00
21) Nitrobenzene-d5	9.44	128	40213	41.62	ng/ul	0.00
24) 2-Nitrophenol-d4	10.16	143	46614	46.10	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.71	165	90821	39.04	ng/ul	0.00
31) 4-Chloroaniline-d4	11.22	131	109772	38.21	ng/ul	0.00
46) Dimethylphthalate-d6	14.27	166	321856	42.35	ng/ul	0.00
49) Acenaphthylene-d8	14.58	160	339235	39.12	ng/ul	0.00
54) 4-Nitrophenol-d4	15.04	143	46348	41.76	ng/ul	0.00
60) Fluorene-d10	15.86	176	283707	41.32	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.96	200	52559	43.84	ng/ul	0.00
73) Anthracene-d10	17.72	188	468573	42.77	ng/ul	0.00
81) Pyrene-d10	19.98	212	543800	41.25	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.07	264	548533	41.91	ng/ul	0.00

Target Compounds Ovalue

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

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