

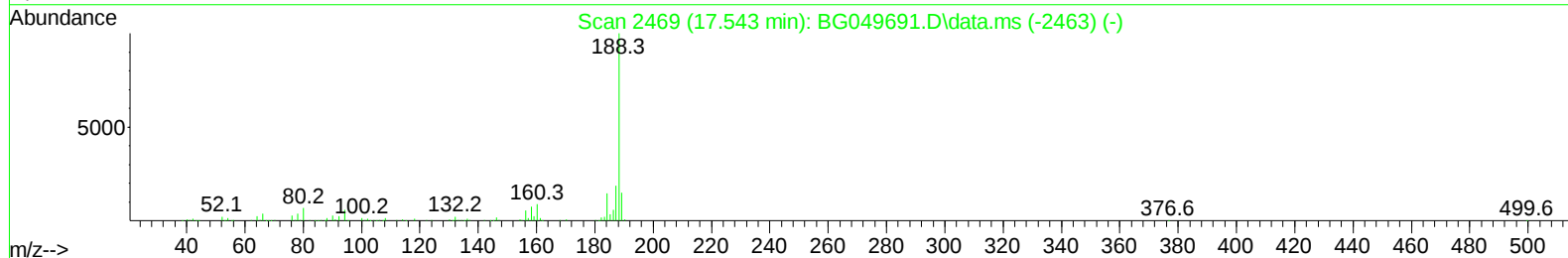
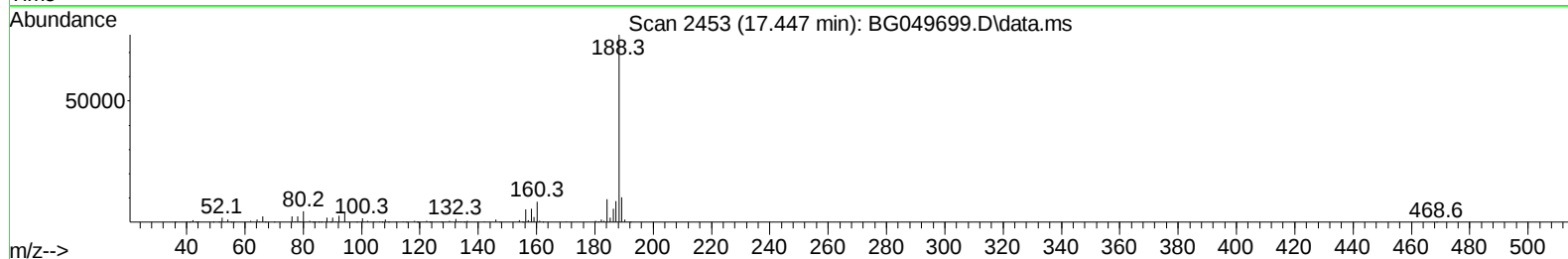
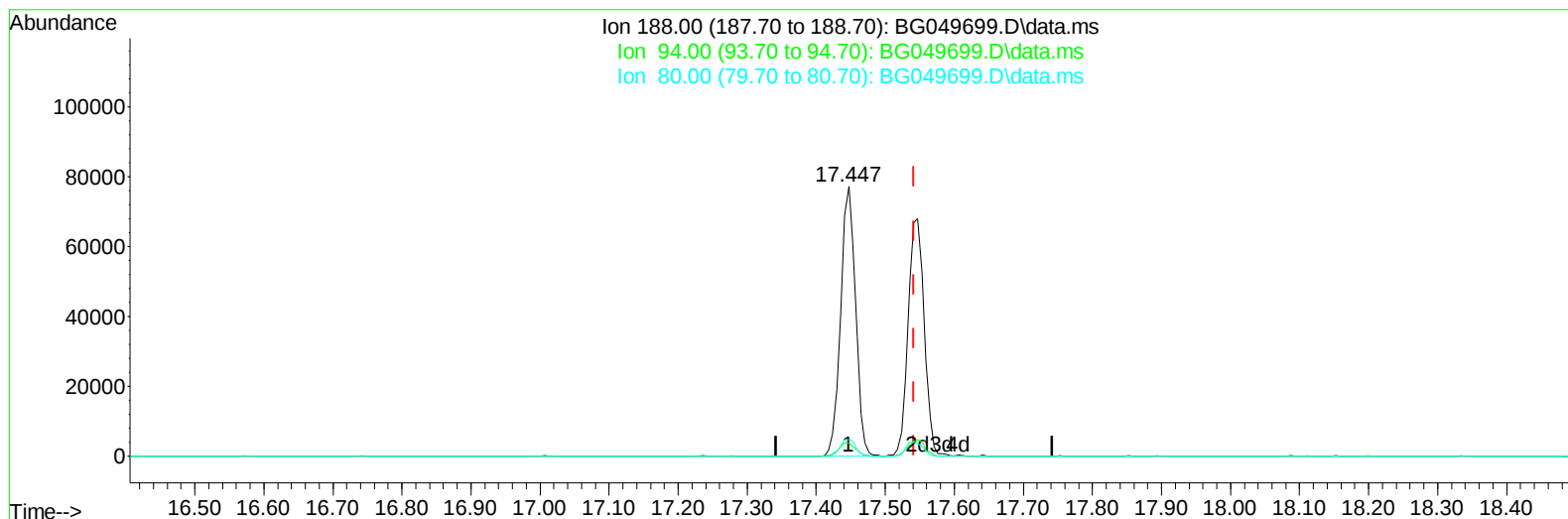
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049699.D
Acq On : 7 Aug 2021 10:37
Operator : CG/JU
Sample : M3164-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JCEN0

Manual Integrations
APPROVED

mohammad
8/9/2021 12:10:29 PM

Quant Time: Aug 07 11:57:01 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 04 16:17:29 2021
Response via : Initial Calibration



TIC: BG049699.D\data.ms

(73) Anthracene-d10 (S)

17.447min (-0.095) 21.18 ng/u1

response 115166

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	4.92
80.00	6.00	6.00
0.00	0.00	0.00

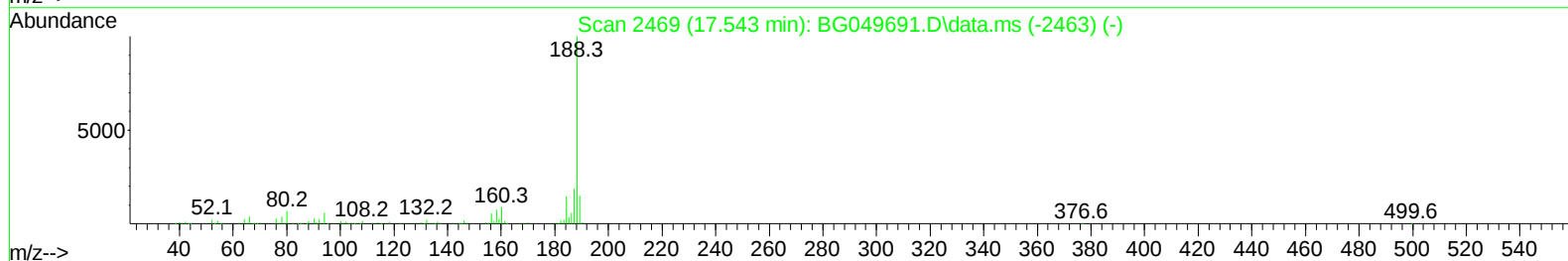
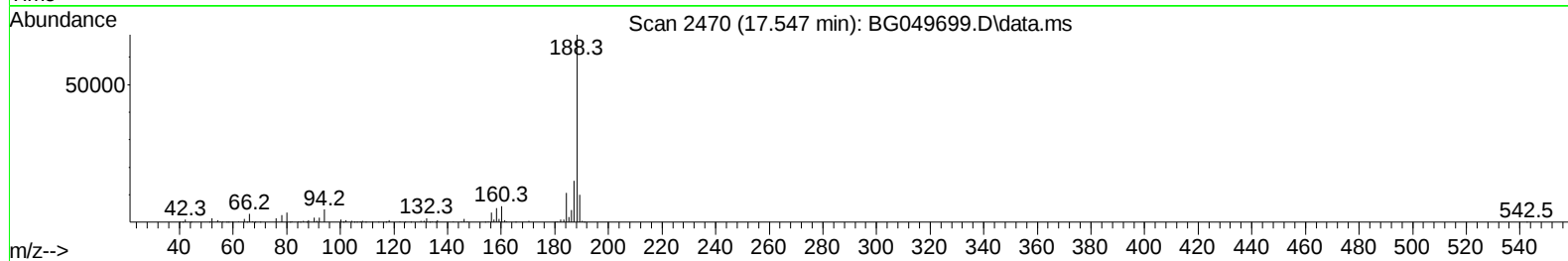
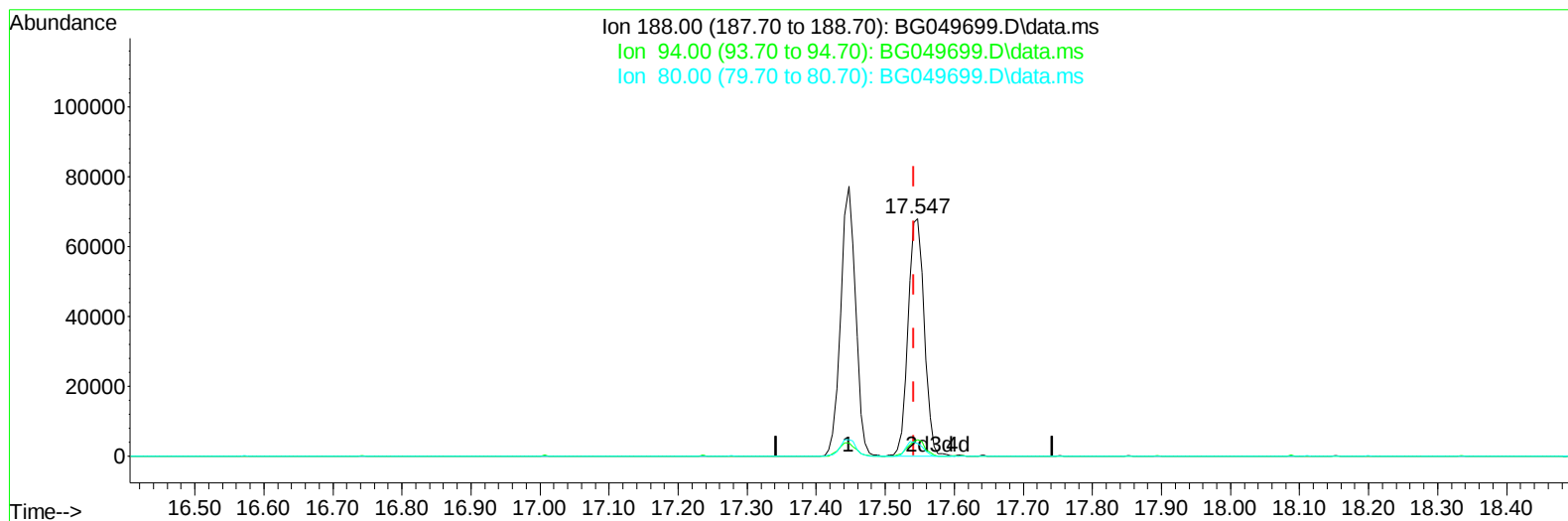
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(73) Anthracene-d10 (S)

17.547min (+ 0.005) 20.25 ng/ul m

response 110064

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	6.87#
80.00	6.00	5.26
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	8.042	152	19568	20.000	ng/ul	0.00
20) Naphthalene-d8	10.868	136	79483	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.692	164	55068	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.447	188	115166	20.000	ng/ul	0.00
79) Chrysene-d12	21.730	240	103063	20.000	ng/ul	0.00
88) Perylene-d12	25.008	264	105798	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.431	96	1266	3.025	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.202	99	27772	15.638	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.361	67	16671	16.343	ng/ul	0.00
11) 2-Chlorophenol-d4	7.572	132	21499	16.914	ng/ul	0.00
15) 4-Methylphenol-d8	8.753	113	17384	12.838	ng/ul	0.00
21) Nitrobenzene-d5	9.217	128	10877	17.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.946	143	12713	17.780	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.486	165	22711	16.845	ng/ul	0.00
31) 4-Chloroaniline-d4	11.003	131	17998	9.751	ng/ul	0.00
46) Dimethylphthalate-d6	14.093	166	79573	18.413	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	98114	19.083	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	8584	12.286	ng/ul	0.02
60) Fluorene-d10	15.691	176	71713	19.305	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.808	200	8403	10.841	ng/ul	0.00
73) Anthracene-d10	17.547	188	110064m	20.246	ng/ul	0.00
81) Pyrene-d10	19.832	212	123669	21.778	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.778	264	102592	17.697	ng/ul	0.00

Target Compounds Qvalue

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

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