

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
 Data File : BM047029.D
 Acq On : 06 Aug 2024 19:13
 Operator : RC/JU
 Sample : P3171-08DL 30X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

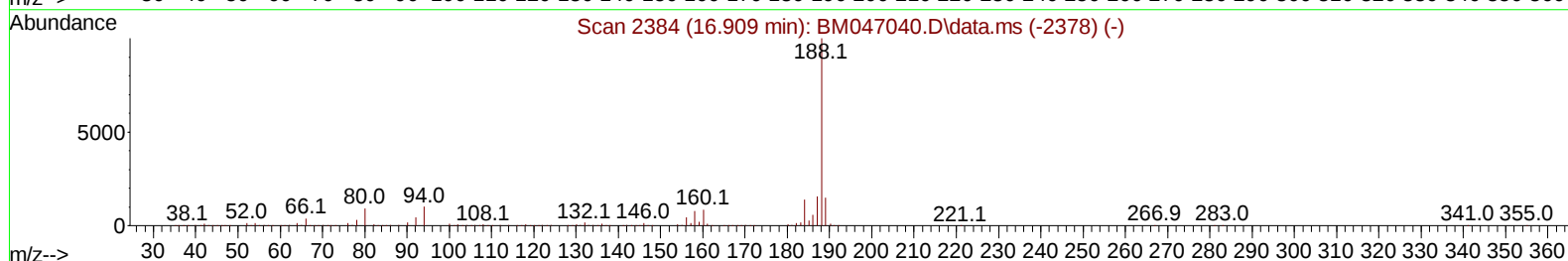
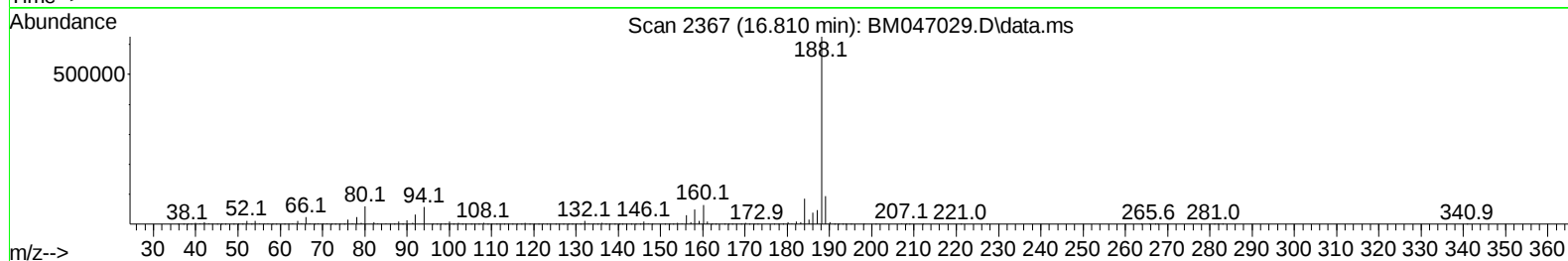
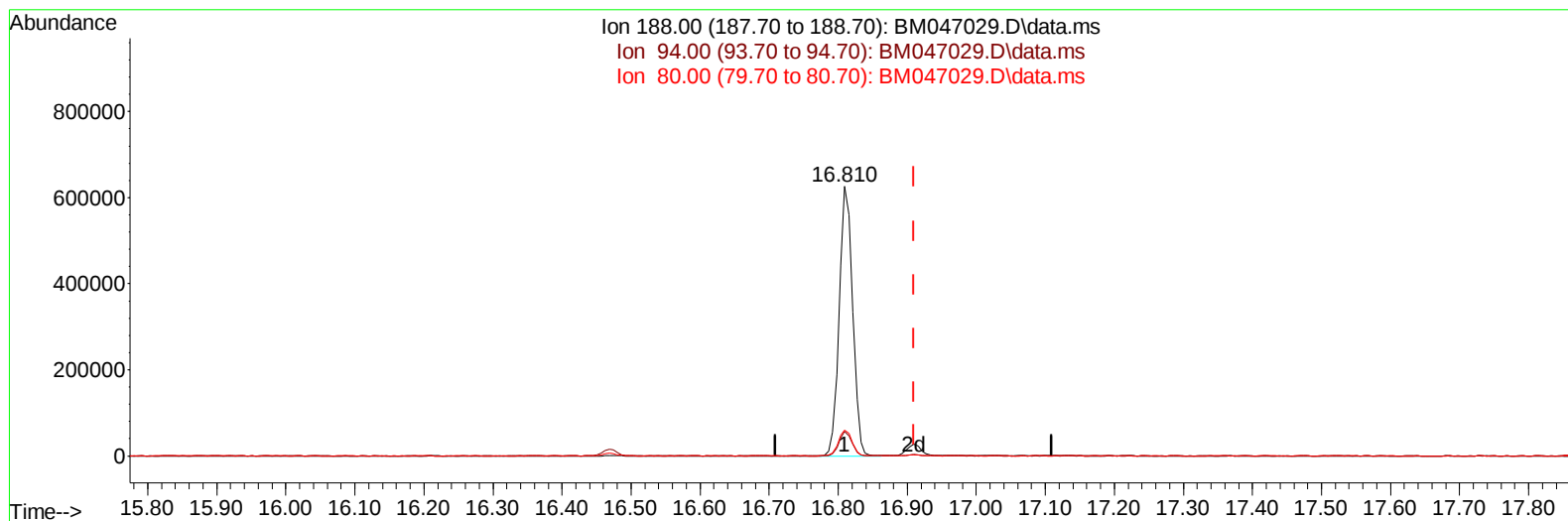
DCVX0DL

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/08/2024

Supervised By :mohammad ahmed 08/30/2024

Quant Time: Aug 07 07:50:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 04:48:25 2024
 Response via : Initial Calibration



TIC: BM047029.D\data.ms

(73) Anthracene-d10 (S)

16.810min (-0.100) 21.09 ng/u1

response 845145

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	9.80	9.09
80.00	8.60	9.53
0.00	0.00	0.00

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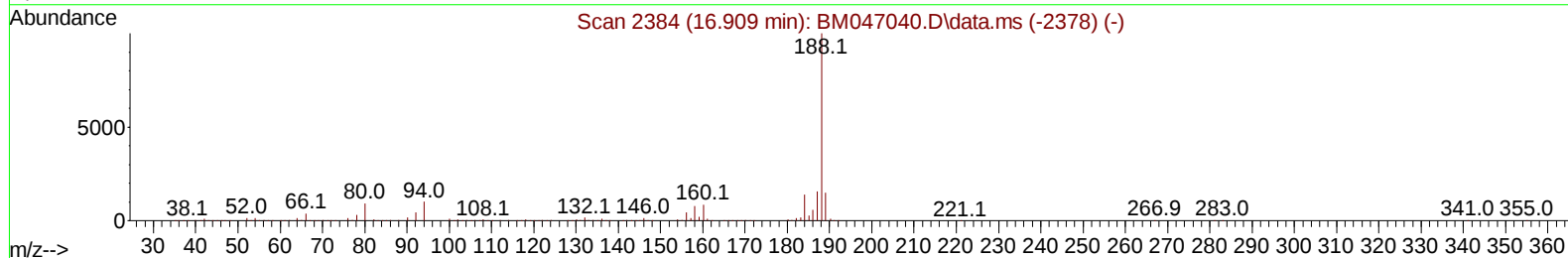
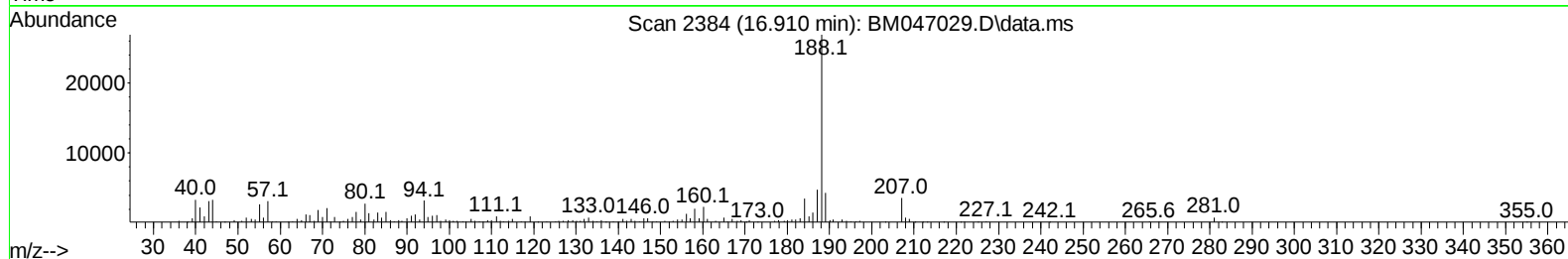
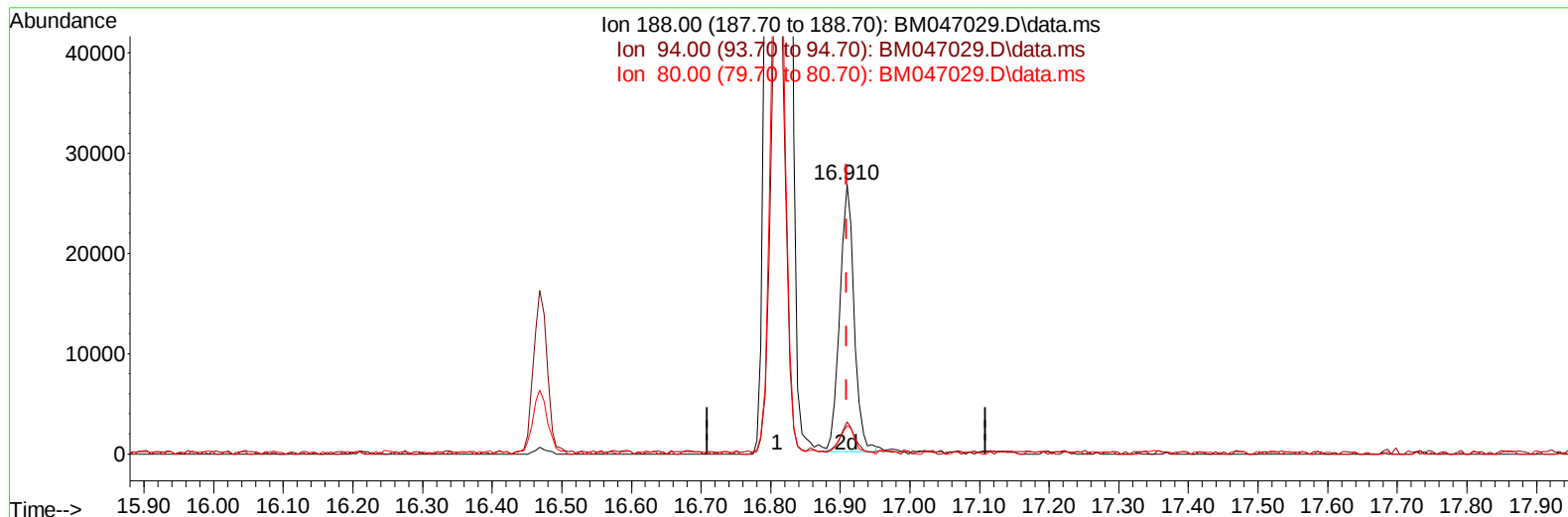
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TIC: BM047029.D\data.ms

(73) Anthracene-d10 (S)

16.910min (+ 0.000) 0.95 ng/u1 m

response 37908

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	9.80	11.88#
80.00	8.60	10.40#
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	7.398	152	161715	20.000	ng/ul	0.00
20) Naphthalene-d8	10.151	136	624527	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.051	164	396787	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.810	188	845145	20.000	ng/ul	0.00
79) Chrysene-d12	21.045	240	764444	20.000	ng/ul	0.00
88) Perylene-d12	23.750	264	880728	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.604	99	10819	0.881	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.757	67	8384	1.037	ng/ul	0.00
11) 2-Chlorophenol-d4	6.940	132	8734	0.839	ng/ul	0.00
15) 4-Methylphenol-d8	8.122	113	7852	0.807	ng/ul	0.00
21) Nitrobenzene-d5	8.545	128	4797	0.957	ng/ul	0.00
24) 2-Nitrophenol-d4	9.251	143	4082	0.746	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.792	165	7672	0.730	ng/ul	0.00
31) 4-Chloroaniline-d4	10.310	131	10324	0.748	ng/ul	0.00
46) Dimethylphthalate-d6	13.480	166	27994	0.923	ng/ul	0.00
49) Acenaphthylene-d8	13.739	160	31663	0.935	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.057	176	25469	0.981	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	16.910	188	37908m	0.946	ng/ul	0.00
81) Pyrene-d10	19.221	212	41425	0.947	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	38649	0.833	ng/ul	-0.01
Target Compounds						
					Qvalue	
36) 2-Methylnaphthalene	11.827	142	26828	1.162	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	14.692	232	65156	8.334	ng/ul	100
71) Pentachlorophenol	16.468	266	913783	125.301	ng/ul	99

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

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