

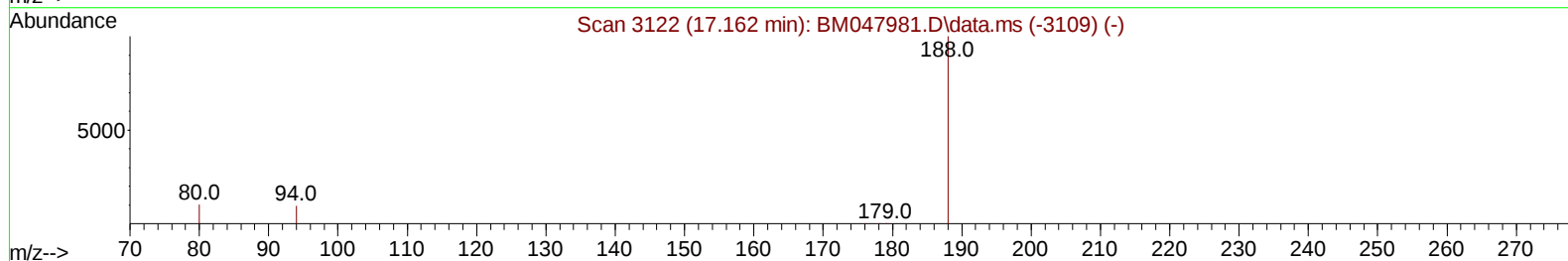
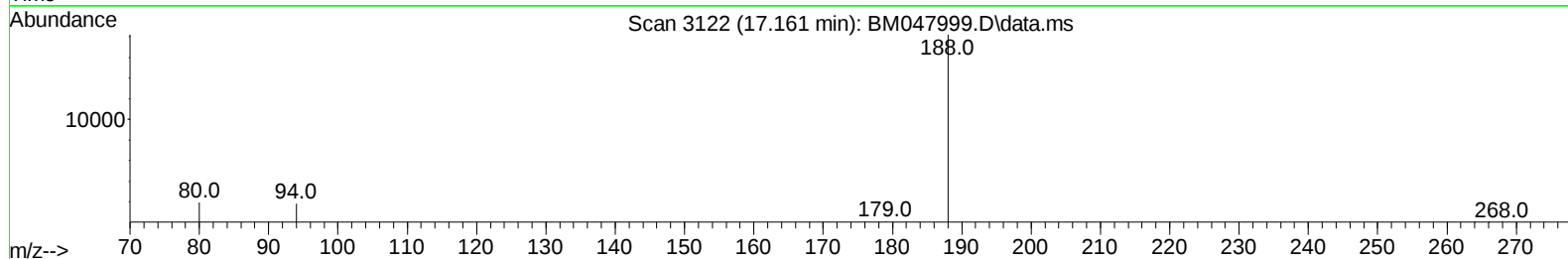
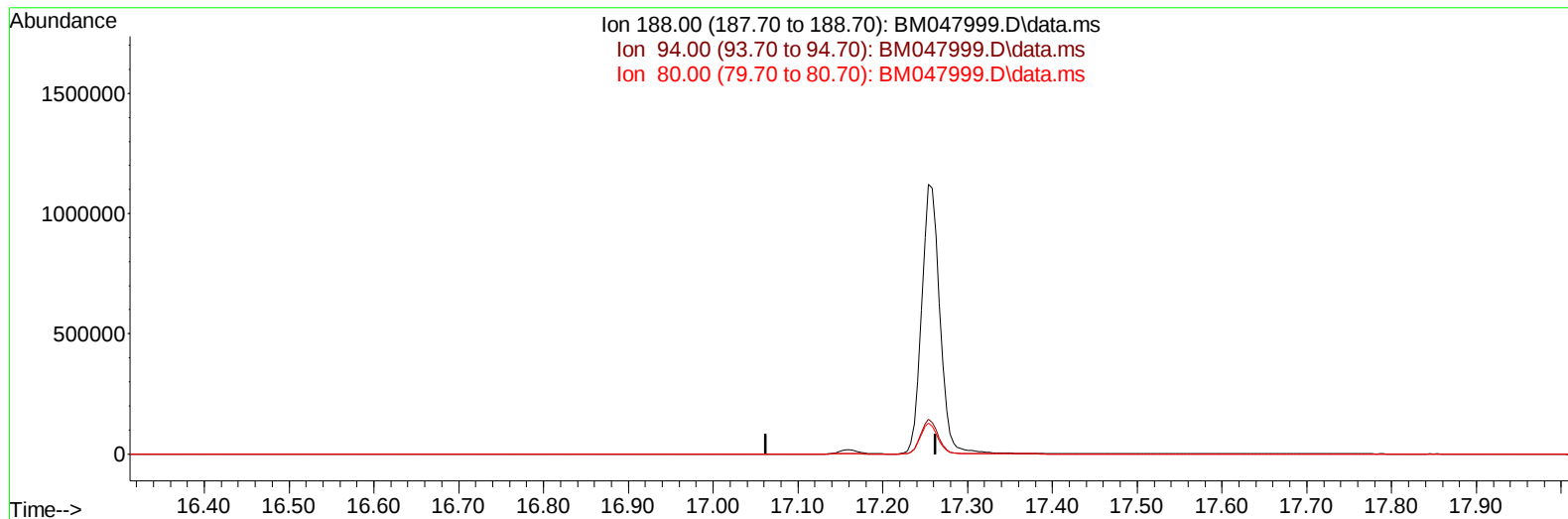
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101024\
Data File : BM047999.D
Acq On : 11 Oct 2024 04:27
Operator : RC/JU
Sample : PB163767BL
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK767

Manual IntegrationsAPPROVED

Quant Time: Oct 11 05:11:10 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 10 18:01:13 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/11/2024
Supervised By :mohammad ahmed 10/14/2024



TIC: BM047999.D\data.ms

(13) Phenanthrene-d10 (I)

17.162min (-17.162) 0.00 ng/u1

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	12.20	0.00#
80.00	13.00	0.00#
0.00	0.00	0.00

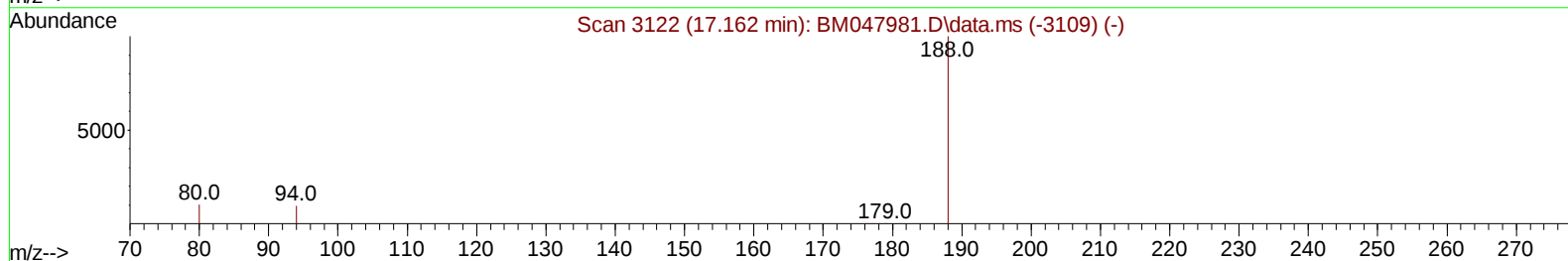
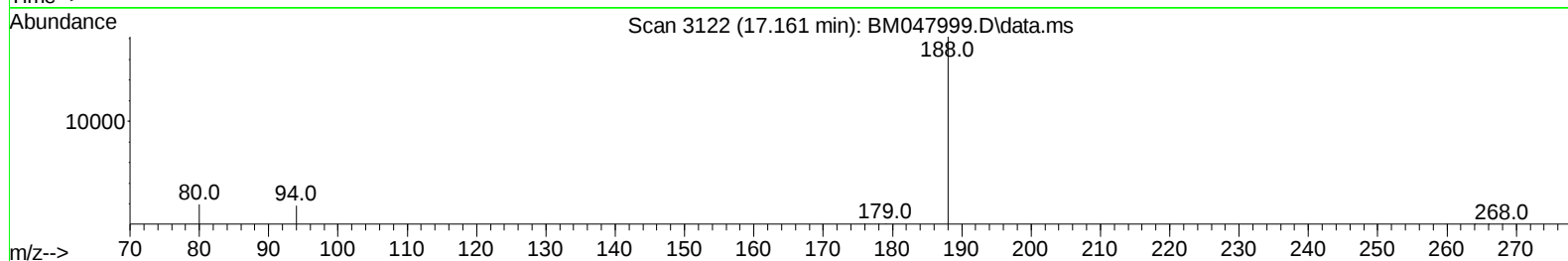
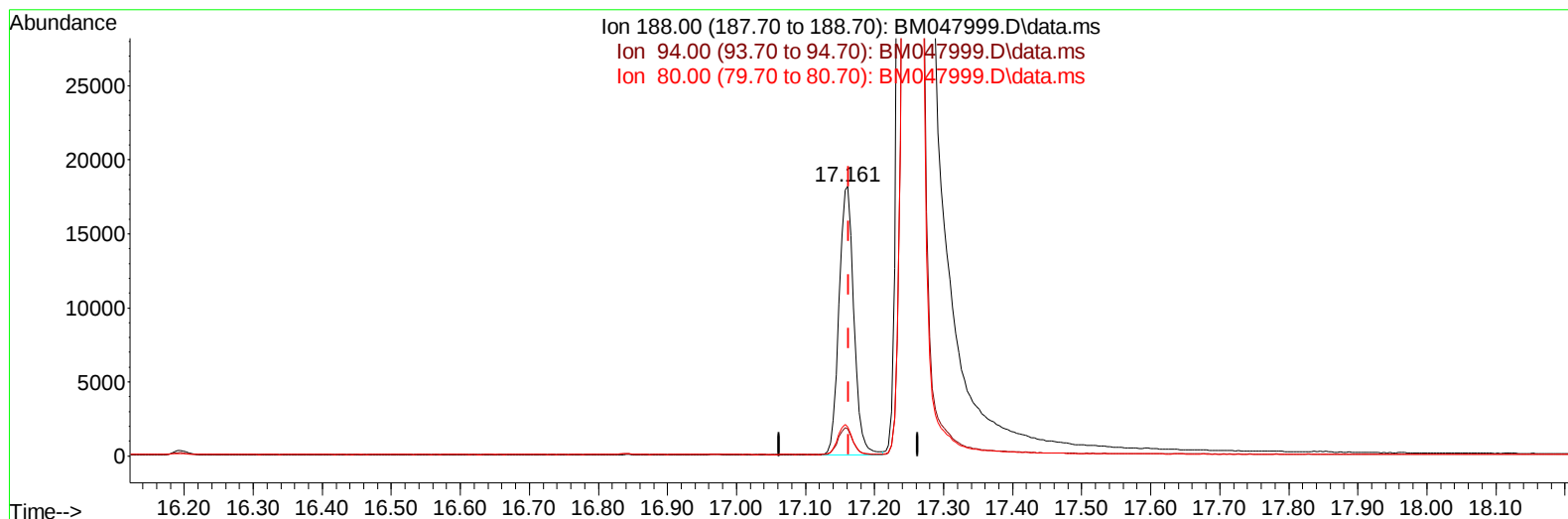
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(13) Phenanthrene-d10 (I)

17.161min (-0.001) 0.40 ng/ul m

response 27238

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	12.20	10.13
80.00	13.00	10.70
0.00	0.00	0.00

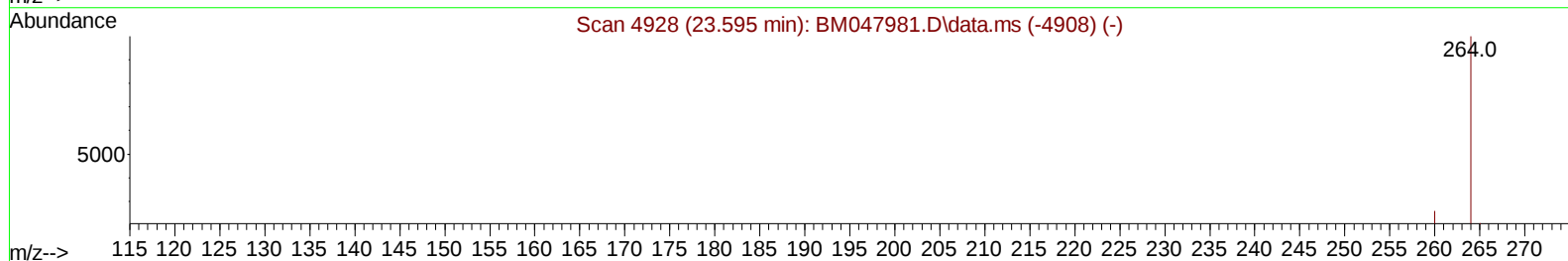
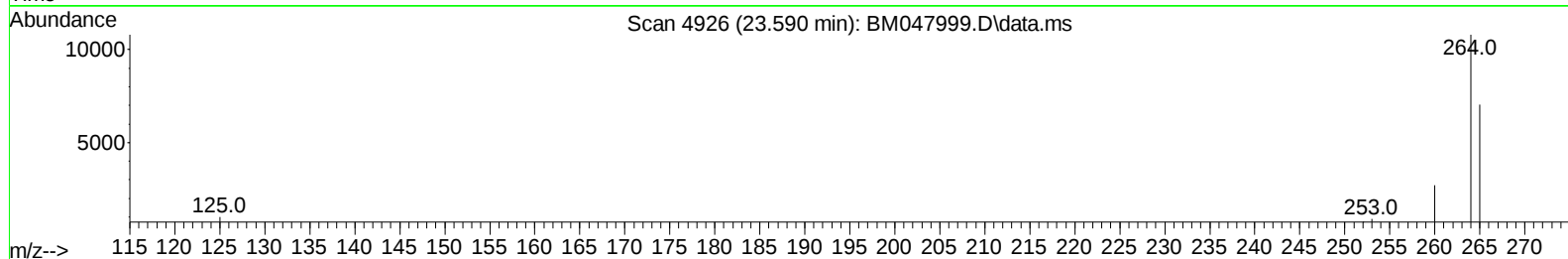
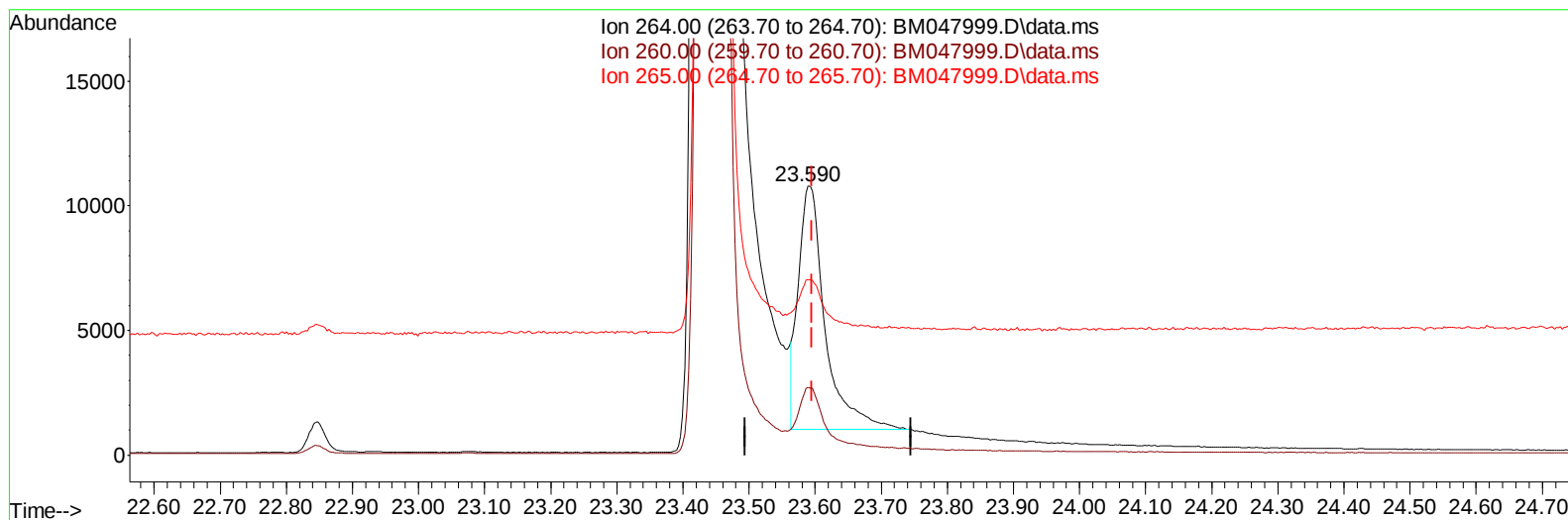
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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TIC: BM047999.D\data.ms

(23) Perylene-d12 (I)

23.590min (-0.005) 0.40 ng/u1

response 27411

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	26.80	25.04
265.00	85.70	65.15#
0.00	0.00	0.00

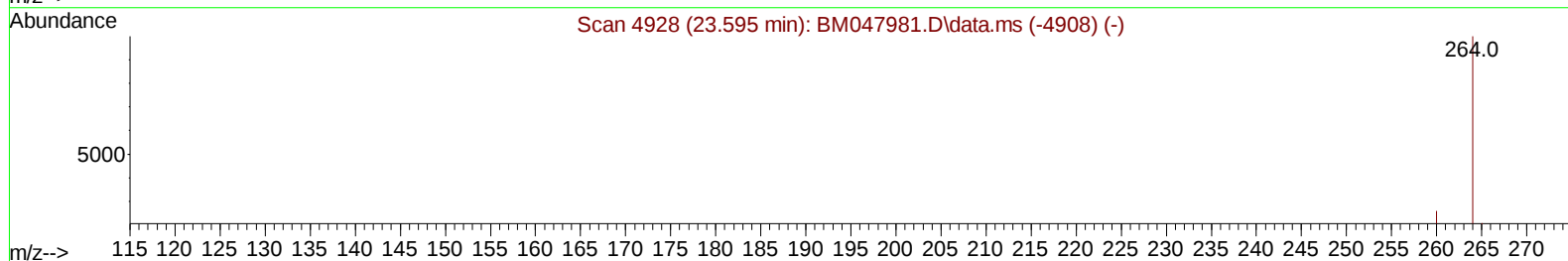
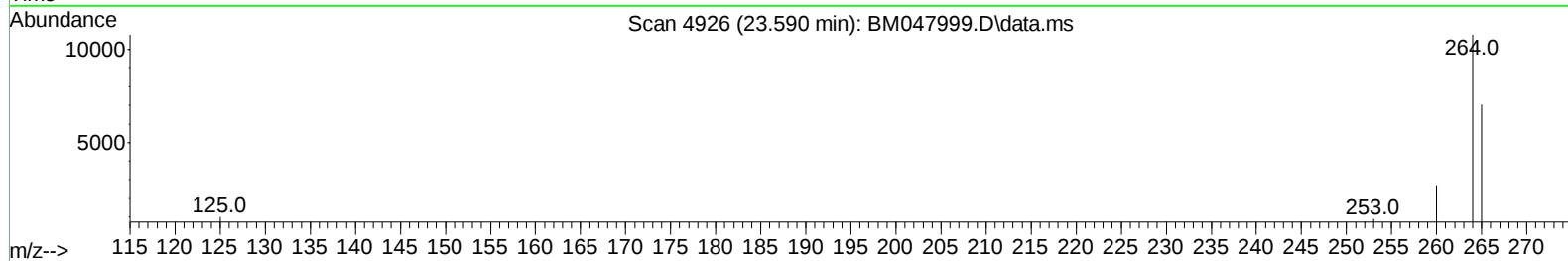
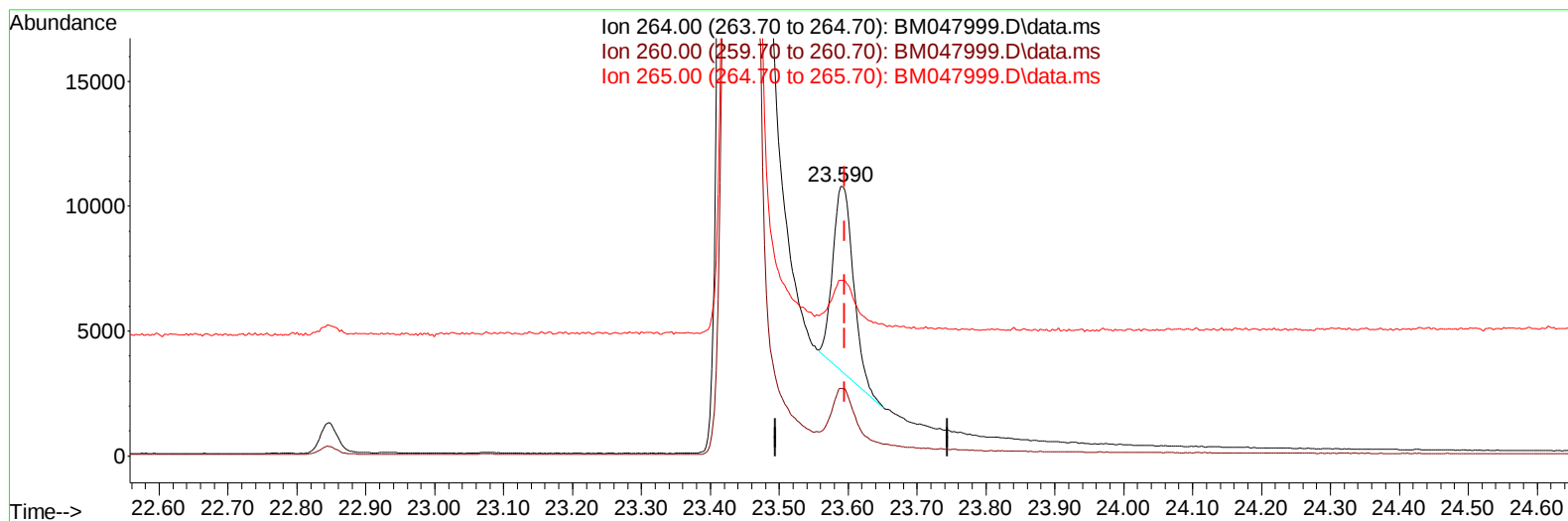
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TIC: BM047999.D\data.ms

(23) Perylene-d12 (I)

23.590min (-0.005) 0.40 ng/ul m

response 15430

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	26.80	25.04
265.00	85.70	65.15#
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	7.792	152	9975	0.400	ng/ul	0.00
4) Naphthalene-d8	10.578	136	29674	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.422	164	14447	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.161	188	27238m	0.400	ng/ul	0.00
17) Chrysene-d12	21.336	240	18881	0.400	ng/ul	0.00
23) Perylene-d12	23.590	264	15430m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	61698	4.814	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.184	152	9172	0.232	ng/ul	0.00
18) Fluoranthene-d10	19.182	212	17537	0.303	ng/ul	0.00

Target Compounds Qvalue

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

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