

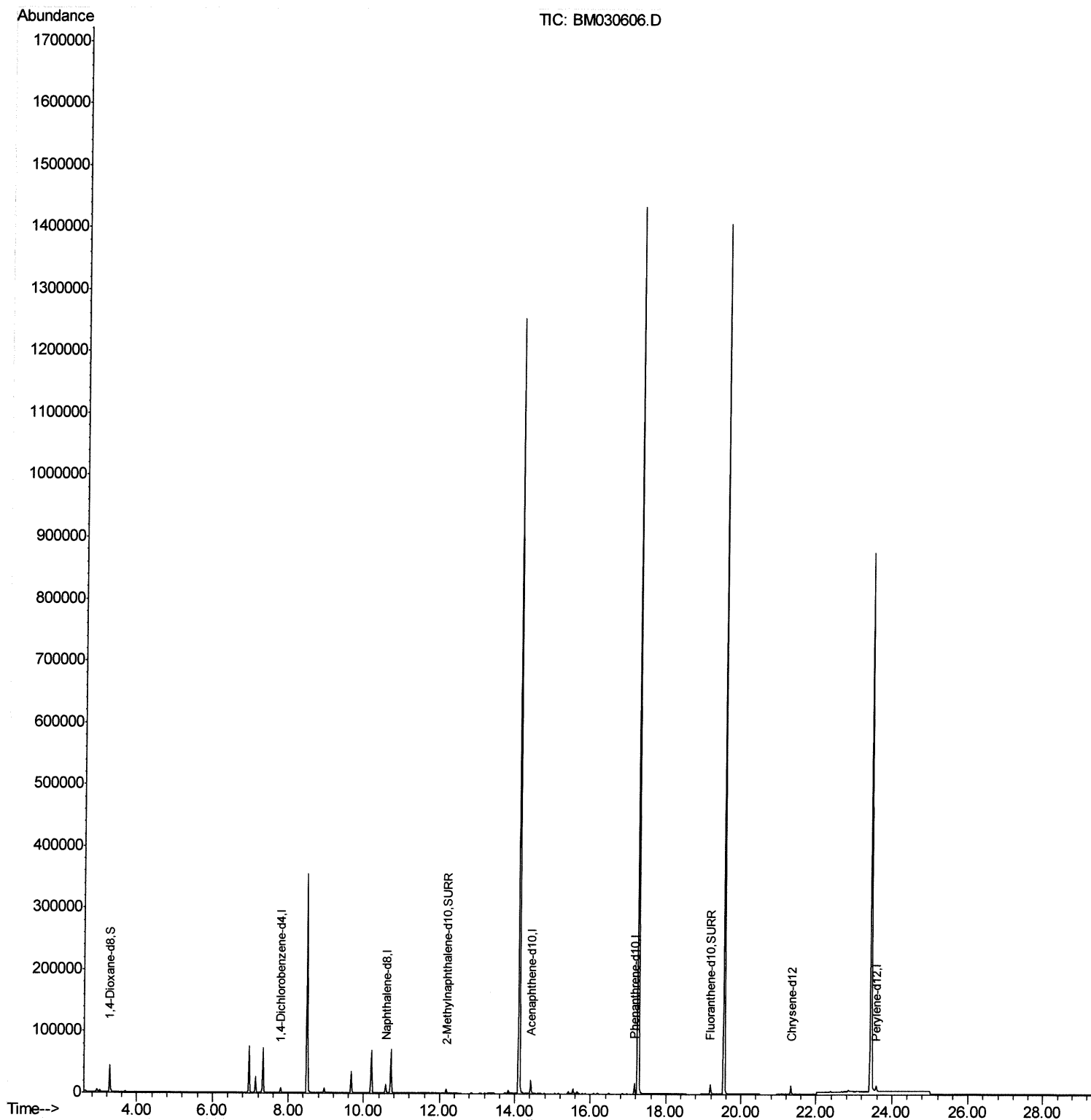
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062521\
Data File : BM030606.D
Acq On : 25 Jun 2021 17:49
Operator : CG/JU
Sample : PB137202BL
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK202

Quant Time: Jun 25 18:21:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 25 12:45:36 2021
Response via : Initial Calibration

Manual Integrations
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mohammad
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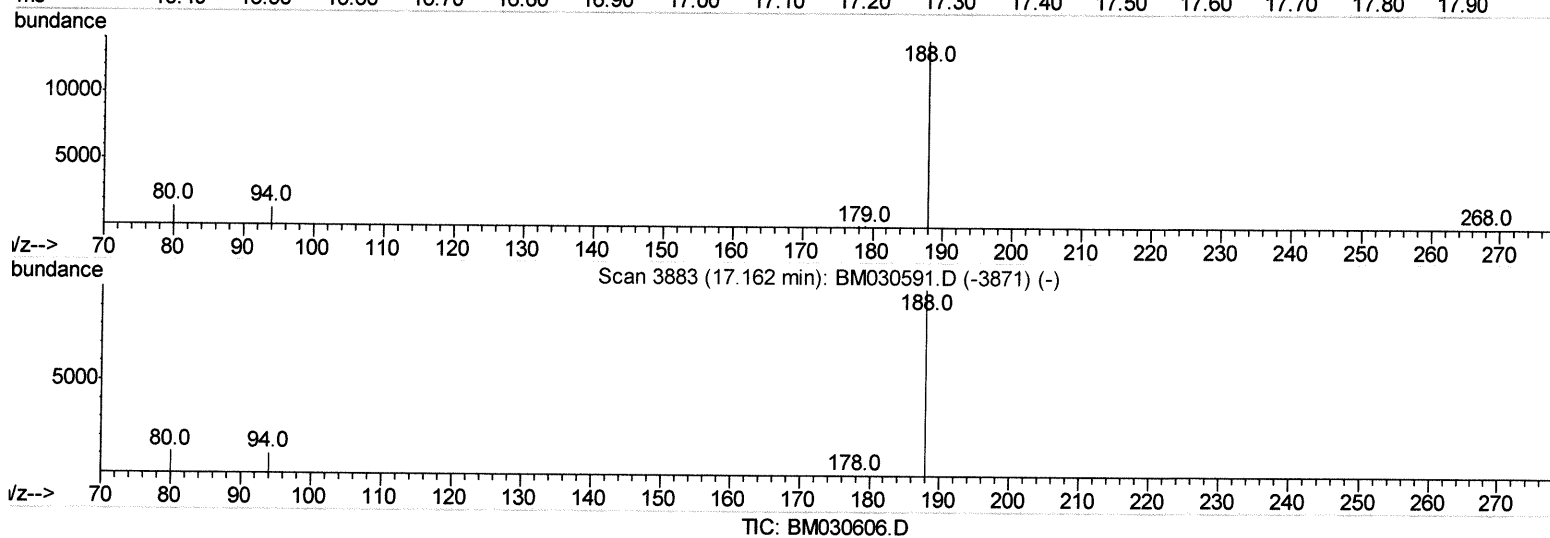
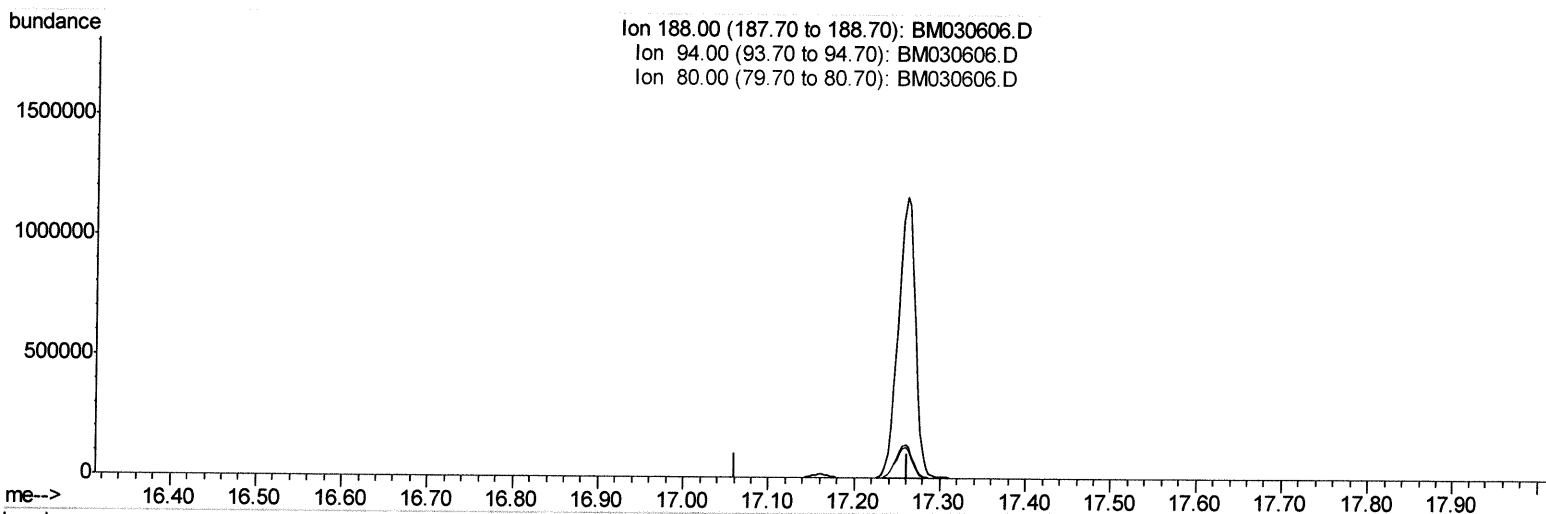
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Quant Time: Jun 25 18:19:01 2021
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(13) Phenanthrene-d10 (I)

17.162min (-17.162) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	10.30	0.00#
80.00	11.60	0.00#
0.00	0.00	0.00

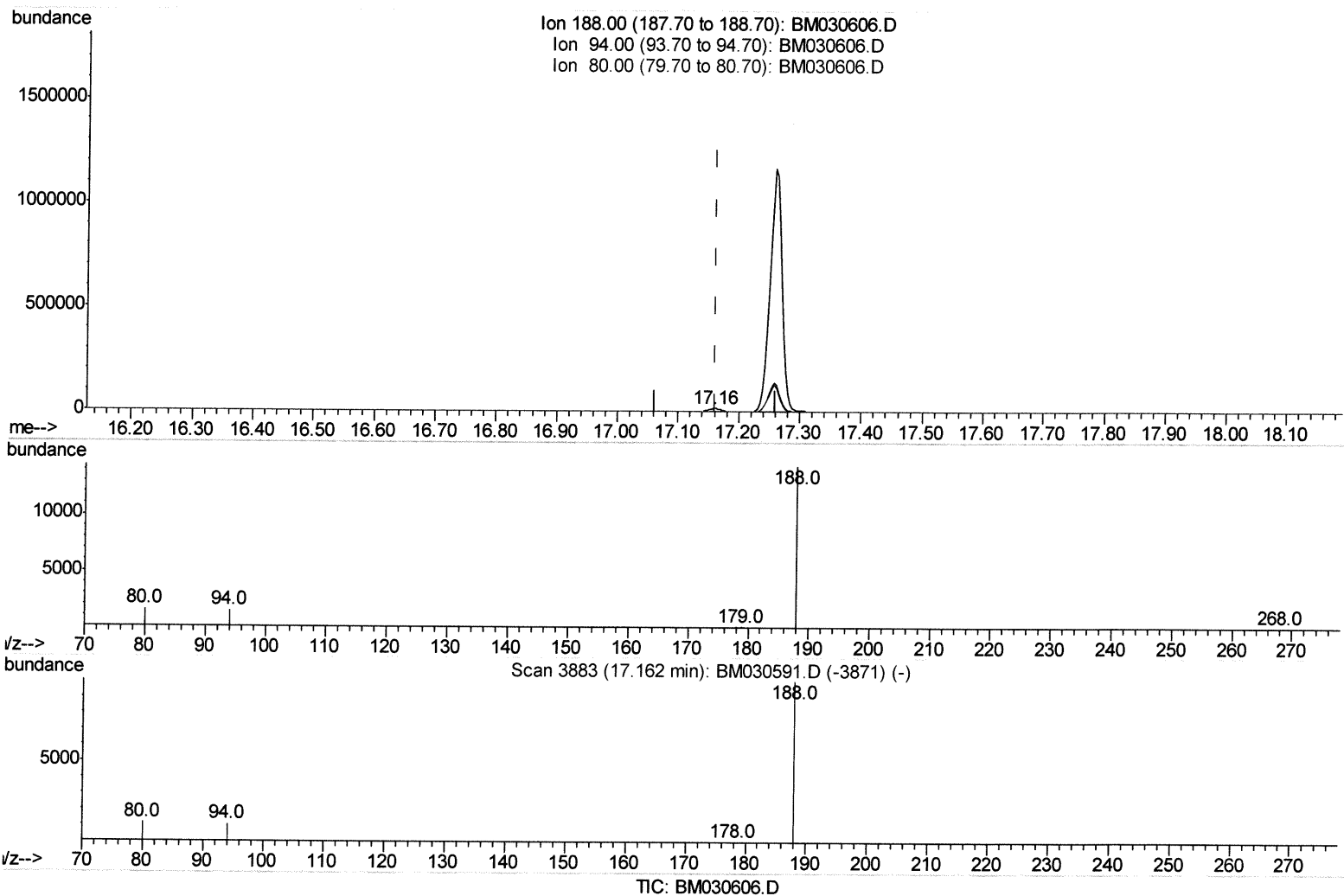
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(13) Phenanthrene-d10 (I)

17.159min (-0.003) 0.40ng/ul m

JU 6/30/21

response 20336

Ion	Exp%	Act%
188.00	100	100
94.00	10.30	10.28
80.00	11.60	10.91
0.00	0.00	0.00

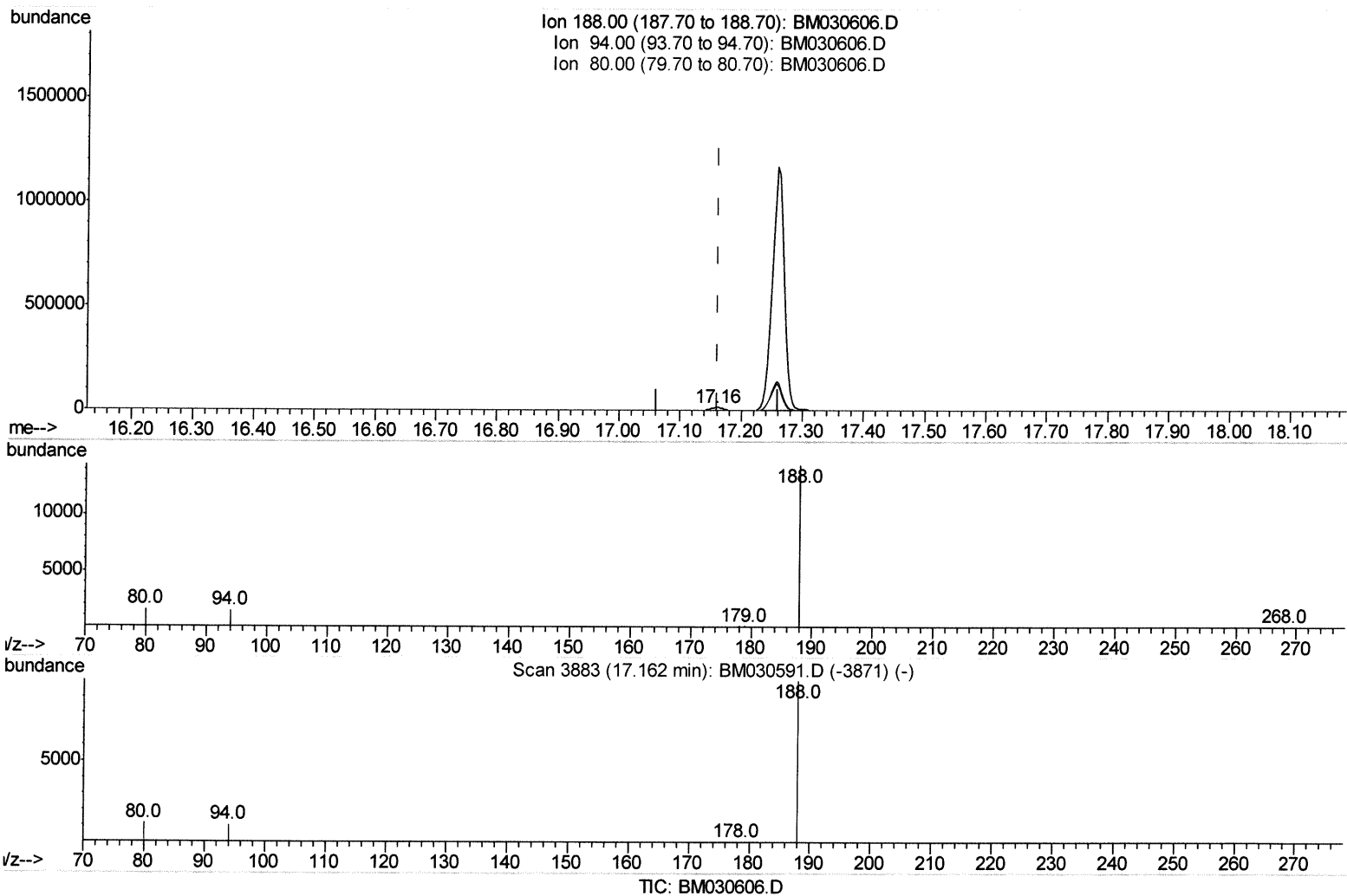
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response 20336

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ju 6/30/21

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	4445	0.40	ng/ul	0.00
4) Naphthalene-d8	10.58	136	18221	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.42	164	11465	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.16	188	20336m	0.40	ng/ul	0.00
17) Chrysene-d12	21.33	240	12262	0.40	ng/ul	0.00
23) Perylene-d12	23.59	264	14003	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.28	96	28546	6.00	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.18	152	10795	0.37	ng/ul	0.00
18) Fluoranthene-d10	19.18	212	18064	0.48	ng/ul	0.00

Target Compounds	Ovalue

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

6/30/21