

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042222\
 Data File : BG053283.D
 Acq On : 23 Apr 2022 5:17
 Operator : CG/JU
 Sample : N2487-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 UST-1-E

Quant Time: Apr 24 01:18:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.106	152	24285	20.000	ng	0.00
21) Naphthalene-d8	10.919	136	96994	20.000	ng	0.00
39) Acenaphthene-d10	14.732	164	77090	20.000	ng	0.00
64) Phenanthrene-d10	17.475	188	199820	20.000	ng	0.00
76) Chrysene-d12	21.758	240	215887	20.000	ng	0.00
86) Perylene-d12	25.053	264	220090	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.674	112	93233	65.810	ng	0.00
7) Phenol-d6	7.266	99	83394	38.172	ng	0.00
23) Nitrobenzene-d5	9.269	82	241203	100.972	ng	0.00
42) 2,4,6-Tribromophenol	16.218	330	181887	148.301	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	545160	97.512	ng	0.00
79) Terphenyl-d14	20.078	244	1153886	90.270	ng	0.00
Target Compounds						
71) Phenanthrene	17.516	178	26643	2.441	ng	Qvalue 93

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

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