

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042222\
 Data File : BG053275.D
 Acq On : 23 Apr 2022 00:08
 Operator : CG/JU
 Sample : N2487-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 UST-1-S

Quant Time: Apr 24 01:17:08 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	24038	20.000	ng	0.00
21) Naphthalene-d8	10.920	136	99534	20.000	ng	# 0.00
39) Acenaphthene-d10	14.732	164	79081	20.000	ng	0.00
64) Phenanthrene-d10	17.475	188	204707	20.000	ng	0.00
76) Chrysene-d12	21.758	240	212132	20.000	ng	0.00
86) Perylene-d12	25.053	264	218058	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.674	112	99581	71.013	ng	0.00
7) Phenol-d6	7.266	99	88919	41.119	ng	0.00
23) Nitrobenzene-d5	9.269	82	248700	101.454	ng	0.00
42) 2,4,6-Tribromophenol	16.218	330	210190	167.063	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	561889	97.974	ng	0.00
79) Terphenyl-d14	20.078	244	1095236	87.198	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.516	178	58093	5.195	ng	97
75) Fluoranthene	19.526	202	34875	2.433	ng	99
78) Pyrene	19.884	202	30221	2.000	ng	97

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

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