

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080222\
 Data File : BG054488.D
 Acq On : 2 Aug 2022 18:14
 Operator : CG/JU
 Sample : N3894-18
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-37

Quant Time: Aug 03 04:14:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.986	152	19014	20.000	ng	0.00
21) Naphthalene-d8	10.794	136	72644	20.000	ng	# 0.00
39) Acenaphthene-d10	14.625	164	53057	20.000	ng	0.00
64) Phenanthrene-d10	17.369	188	129788	20.000	ng	# 0.00
76) Chrysene-d12	21.640	240	145853	20.000	ng	# 0.00
86) Perylene-d12	24.848	264	150478	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.571	112	104706	114.975	ng	0.00
7) Phenol-d6	7.187	99	140887	112.903	ng	0.01
23) Nitrobenzene-d5	9.149	82	96212	72.125	ng	0.00
42) 2,4,6-Tribromophenol	16.123	330	127237	114.216	ng	0.00
45) 2-Fluorobiphenyl	13.244	172	293453	78.024	ng	0.00
79) Terphenyl-d14	19.983	244	579103	78.766	ng	0.00
Target Compounds						
78) Pyrene	19.789	202	22976	2.745	ng	Qvalue 98
88) Benzo(b)fluoranthene	23.838	252	18446	2.073	ng	# 93

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

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