

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072922\
 Data File : BG054447.D
 Acq On : 29 Jul 2022 13:24
 Operator : CG/JU
 Sample : N3895-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-43

Quant Time: Jul 30 03:17:55 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	7.982	152	16647	20.000	ng	# 0.00
21) Naphthalene-d8	10.797	136	66234	20.000	ng	# 0.00
39) Acenaphthene-d10	14.621	164	51374	20.000	ng	0.00
64) Phenanthrene-d10	17.371	188	126715	20.000	ng	# 0.00
76) Chrysene-d12	21.637	240	142372	20.000	ng	# 0.00
86) Perylene-d12	24.839	264	151594	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	83406	104.609	ng	0.00
7) Phenol-d6	7.177	99	115563	105.777	ng	0.00
23) Nitrobenzene-d5	9.151	82	78824	64.808	ng	0.00
42) 2,4,6-Tribromophenol	16.114	330	113159	104.907	ng	0.00
45) 2-Fluorobiphenyl	13.241	172	251346	69.018	ng	0.00
79) Terphenyl-d14	19.980	244	510572	71.143	ng	0.00
Target Compounds						
78) Pyrene	19.786	202	20232	2.477	ng	Qvalue 95
84) Bis(2-ethylhexyl)phtha...	21.508	149	8097	2.248	ng	# 94
88) Benzo(b)fluoranthene	23.828	252	21139	2.358	ng	# 96

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

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