

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080122\
 Data File : BG054460.D
 Acq On : 1 Aug 2022 11:46
 Operator : CG/JU
 Sample : PB146613BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB146613BL

Quant Time: Aug 02 05:45:07 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	17227	20.000	ng	# 0.00
21) Naphthalene-d8	10.796	136	62077	20.000	ng	# 0.00
39) Acenaphthene-d10	14.621	164	49838	20.000	ng	0.00
64) Phenanthrene-d10	17.371	188	133550	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	160100	20.000	ng	# 0.00
86) Perylene-d12	24.833	264	170912	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	117285	142.147	ng	0.00
7) Phenol-d6	7.177	99	151277	133.804	ng	0.00
23) Nitrobenzene-d5	9.151	82	104051	91.279	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	132742	126.854	ng	0.00
45) 2-Fluorobiphenyl	13.240	172	312155	88.357	ng	0.00
79) Terphenyl-d14	19.980	244	768010	95.164	ng	0.00

Target Compounds	Qvalue

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

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