

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060814.D
 Acq On : 6 Apr 2024 16:54
 Operator : MA/JU
 Sample : P1996-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP3

Quant Time: Apr 08 02:08:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 03:12:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.953	152	67792	20.000	ng	0.00
21) Naphthalene-d8	10.750	136	256132	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	187473	20.000	ng	0.00
64) Phenanthrene-d10	17.325	188	437424	20.000	ng	-0.01
76) Chrysene-d12	21.590	240	434925	20.000	ng	-0.01
86) Perylene-d12	24.728	264	502746	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.544	112	494075	121.412	ng	0.00
7) Phenol-d6	7.119	99	636353	105.346	ng	0.00
23) Nitrobenzene-d5	9.105	82	481820	107.341	ng	0.00
42) 2,4,6-Tribromophenol	16.067	330	376871	177.091	ng	0.00
45) 2-Fluorobiphenyl	13.206	172	1146872	89.783	ng	-0.01
79) Terphenyl-d14	19.957	244	2169273	87.825	ng	0.00

Target Compounds	Qvalue

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

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