

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040424\
 Data File : BG060783.D
 Acq On : 4 Apr 2024 18:15
 Operator : MA/JU
 Sample : P1957-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM-DRUMS-040224

Quant Time: Apr 04 19:11:48 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	7.956	152	55280	20.000	ng	0.00
21) Naphthalene-d8	10.752	136	205806	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	156224	20.000	ng	0.00
64) Phenanthrene-d10	17.333	188	399000	20.000	ng	0.00
76) Chrysene-d12	21.593	240	434058	20.000	ng	-0.01
86) Perylene-d12	24.736	264	518686	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.541	112	389021	117.233	ng	0.00
7) Phenol-d6	7.116	99	450399	91.439	ng	0.00
23) Nitrobenzene-d5	9.107	82	348019	96.491	ng	0.00
42) 2,4,6-Tribromophenol	16.070	330	287823	162.301	ng	0.00
45) 2-Fluorobiphenyl	13.208	172	817291	76.780	ng	-0.01
79) Terphenyl-d14	19.959	244	1802074	73.104	ng	0.00

Target Compounds	Qvalue

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

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