

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040124\
 Data File : BG060719.D
 Acq On : 1 Apr 2024 17:31
 Operator : MA/JU
 Sample : P1894-04DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RB40858RTDL

Quant Time: Apr 01 18:17:13 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	7.955	152	65999	20.000	ng	0.00
21) Naphthalene-d8	10.752	136	314706	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	235613	20.000	ng	0.00
64) Phenanthrene-d10	17.333	188	492142	20.000	ng	# 0.00
76) Chrysene-d12	21.592	240	458733	20.000	ng	#-0.01
86) Perylene-d12	24.730	264	542151	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.535	112	249443	62.962	ng	0.00
7) Phenol-d6	7.109	99	319839	54.387	ng	0.00
23) Nitrobenzene-d5	9.107	82	173586	31.474	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	140588	52.564	ng	0.00
45) 2-Fluorobiphenyl	13.208	172	538766	33.560	ng	-0.01
79) Terphenyl-d14	19.959	244	855014	32.819	ng	0.00
Target Compounds						
70) Pentachlorophenol	16.974	266	112034	30.453	ng	99
84) Bis(2-ethylhexyl)phtha...	21.528	149	48918	2.622	ng	# 95

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

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