

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060807.D
 Acq On : 6 Apr 2024 12:10
 Operator : MA/JU
 Sample : P1995-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 280648

Quant Time: Apr 08 02:03:21 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	7.954	152	52946	20.000	ng	0.00
21) Naphthalene-d8	10.751	136	200176	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	156566	20.000	ng	0.00
64) Phenanthrene-d10	17.331	188	386856	20.000	ng	0.00
76) Chrysene-d12	21.591	240	385156	20.000	ng	-0.01
86) Perylene-d12	24.734	264	439858	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.539	112	369848	116.369	ng	0.00
7) Phenol-d6	7.114	99	460885	97.692	ng	0.00
23) Nitrobenzene-d5	9.105	82	351832	100.292	ng	0.00
42) 2,4,6-Tribromophenol	16.068	330	320092	180.102	ng	0.00
45) 2-Fluorobiphenyl	13.206	172	881776	82.657	ng	-0.01
79) Terphenyl-d14	19.957	244	1897302	86.740	ng	0.00

Target Compounds	Qvalue

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

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