

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040524\
 Data File : BG060802.D
 Acq On : 5 Apr 2024 19:06
 Operator : MA/JU
 Sample : P1995-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MOO-23-00119

Quant Time: Apr 05 23:10:06 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	80100	20.000	ng	0.00
21) Naphthalene-d8	10.751	136	299925	20.000	ng	0.00
39) Acenaphthene-d10	14.723	164	2108759	20.000	ng	# 0.13
64) Phenanthrene-d10	17.397	188	283373	20.000	ng	# 0.06
76) Chrysene-d12	21.639	240	284093	20.000	ng	# 0.03
86) Perylene-d12	24.764	264	386726	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	329276	68.482	ng	0.00
7) Phenol-d6	7.115	99	464258	65.047	ng	0.00
23) Nitrobenzene-d5	9.106	82	306615	58.334	ng	0.00
42) 2,4,6-Tribromophenol	16.122	330	166031	6.936	ng	0.05
45) 2-Fluorobiphenyl	13.213	172	676482	4.708	ng	0.00
79) Terphenyl-d14	20.011	244	865175	53.624	ng	0.05
Target Compounds						Qvalue
10) Phenol	7.144	94	14768	2.027	ng	95
22) Acetophenone	8.771	105	29796	3.603	ng	# 89
37) 2-Methylnaphthalene	12.408	142	27058	2.268	ng	# 82
38) 1-Methylnaphthalene	12.632	142	24943	2.214	ng	# 98
84) Bis(2-ethylhexyl)phtha...	21.545	149	129122	11.177	ng	# 75

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

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