

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072823\
 Data File : BG058528.D
 Acq On : 28 Jul 2023 21:45
 Operator : MA/JU
 Sample : 03719-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B-P22B

Quant Time: Jul 29 01:05:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 22:29:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	51167	20.000	ng	0.00
21) Naphthalene-d8	10.614	136	213956	20.000	ng	0.00
39) Acenaphthene-d10	14.462	164	143863	20.000	ng	0.00
64) Phenanthrene-d10	17.206	188	319370	20.000	ng	0.00
76) Chrysene-d12	21.448	240	281478	20.000	ng	0.00
86) Perylene-d12	24.515	264	351762	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.385	112	274398	81.968	ng	0.00
7) Phenol-d6	6.983	99	382972	82.215	ng	0.00
23) Nitrobenzene-d5	8.975	82	216671	49.749	ng	0.00
42) 2,4,6-Tribromophenol	15.949	330	166629	70.665	ng	0.00
45) 2-Fluorobiphenyl	13.082	172	457217	47.627	ng	0.00
79) Terphenyl-d14	19.821	244	727330	48.592	ng	0.00
Target Compounds						
						Qvalue
50) Dimethylphthalate	13.910	163	71130	6.583	ng	100
71) Phenanthrene	17.247	178	38063	2.513	ng	98
75) Fluoranthene	19.263	202	65993	3.627	ng	98
78) Pyrene	19.627	202	128210	6.693	ng	96
81) Benzo(a)anthracene	21.425	228	61655	3.183	ng	98
83) Chrysene	21.490	228	60961	3.282	ng	99
88) Benzo(b)fluoranthene	23.540	252	54184	2.840	ng	# 95
90) Benzo(a)pyrene	24.369	252	52609	2.808	ng	97

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

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