

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111721\
 Data File : BG051096.D
 Acq On : 17 Nov 2021 17:41
 Operator : CG/JU
 Sample : M4673-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RO-72-11933

Quant Time: Nov 18 00:03:42 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 03:55:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.223	152	41431	20.000	ng	0.00
21) Naphthalene-d8	11.050	136	168766	20.000	ng	# 0.00
39) Acenaphthene-d10	14.863	164	93134	20.000	ng	0.00
64) Phenanthrene-d10	17.648	188	158567	20.000	ng	# 0.05
76) Chrysene-d12	21.937	240	153704	20.000	ng	# 0.04
86) Perylene-d12	25.321	264	163536	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.762	112	224244	80.965	ng	0.00
7) Phenol-d6	7.377	99	304011	77.768	ng	0.00
23) Nitrobenzene-d5	9.399	82	222140	58.266	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	89311	95.825	ng	0.02
45) 2-Fluorobiphenyl	13.476	172	379016	64.301	ng	0.00
79) Terphenyl-d14	20.239	244	507473	60.413	ng	0.05
Target Compounds						
						Qvalue
58) Fluorene	15.920	166	16806	2.519	ng	# 1
71) Phenanthrene	17.689	178	42294	5.000	ng	# 1
74) Di-n-butylphthalate	18.611	149	105157	10.654	ng	# 50
75) Fluoranthene	19.710	202	65453	6.308	ng	# 19
78) Pyrene	20.068	202	70380	6.104	ng	# 46
81) Benzo(a)anthracene	21.913	228	35461	3.360	ng	# 82
83) Chrysene	21.984	228	34678	3.483	ng	# 79
84) Bis(2-ethylhexyl)phtha...	21.755	149	18145	2.600	ng	# 56
87) Indeno(1,2,3-cd)pyrene	29.234	276	34926	3.009	ng	# 92
88) Benzo(b)fluoranthene	24.234	252	55707	5.412	ng	# 85
90) Benzo(a)pyrene	25.162	252	42930	4.885	ng	# 91
92) Benzo(g,h,i)perylene	30.468	276	31611	3.361	ng	# 91

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

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