

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062822\
 Data File : BG054117.D
 Acq On : 28 Jun 2022 14:03
 Operator : CG/JU
 Sample : N3472-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20071

Quant Time: Jun 28 14:39:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.070	152	32909	20.000	ng	0.00
21) Naphthalene-d8	10.878	136	124511	20.000	ng	# 0.00
39) Acenaphthene-d10	14.733	164	68315	20.000	ng	# 0.03
64) Phenanthrene-d10	17.524	188	124783	20.000	ng	# 0.08
76) Chrysene-d12	21.801	240	149304	20.000	ng	# 0.08
86) Perylene-d12	25.074	264	150377	20.000	ng	0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	99109	57.579	ng	0.00
7) Phenol-d6	7.230	99	130330	55.815	ng	0.00
23) Nitrobenzene-d5	9.227	82	88820	40.088	ng	0.00
42) 2,4,6-Tribromophenol	16.249	330	54378	49.543	ng	0.06
45) 2-Fluorobiphenyl	13.328	172	194817	41.468	ng	0.00
79) Terphenyl-d14	20.126	244	244937	33.042	ng	0.08
Target Compounds						
						Qvalue
46) 1,1'-Biphenyl	13.540	154	41909	8.807	ng	93
71) Phenanthrene	17.571	178	94333	14.714	ng	# 1
75) Fluoranthene	19.586	202	32653	3.880	ng	# 1
78) Pyrene	19.950	202	44192	4.901	ng	# 1
81) Benzo(a)anthracene	21.783	228	27534	2.848	ng	# 43
83) Chrysene	21.848	228	27106	2.942	ng	# 47
84) Bis(2-ethylhexyl)phtha...	21.672	149	44995	10.884	ng	# 57
88) Benzo(b)fluoranthene	24.034	252	36514	3.994	ng	# 1
90) Benzo(a)pyrene	24.927	252	21755	2.820	ng	# 1

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

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