

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020623\
 Data File : BG056562.D
 Acq On : 6 Feb 2023 19:49
 Operator : CG/JU
 Sample : 01418-11DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JPP-20-020223DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/07/2023
 Supervised By : Jagrut Upadhyay 02/07/2023

Quant Time: Feb 07 04:35:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 07 04:30:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.278	152	49170	20.000	ng	0.00
21) Naphthalene-d8	11.110	136	179320	20.000	ng	0.00
39) Acenaphthene-d10	14.894	164	131334	20.000	ng	0.00
64) Phenanthrene-d10	17.637	188	291717	20.000	ng	0.00
76) Chrysene-d12	21.927	240	264804	20.000	ng	0.00
86) Perylene-d12	25.358	264	292454	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.804	112	42606	15.940	ng	0.01
7) Phenol-d6	7.426	99	60106	15.177	ng	0.00
23) Nitrobenzene-d5	9.453	82	32448	10.275	ng	0.00
42) 2,4,6-Tribromophenol	16.380	330	25348	14.518	ng	0.00
45) 2-Fluorobiphenyl	13.519	172	101716	10.738	ng	0.00
79) Terphenyl-d14	20.211	244	154310	10.235	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.679	178	295564	19.358	ng	98
72) Anthracene	17.767	178	68082	4.344	ng	100
75) Fluoranthene	19.670	202	515262	26.739	ng	99
78) Pyrene	20.035	202	570495	30.292	ng	99
81) Benzo(a)anthracene	21.909	228	276684	14.946	ng	99
83) Chrysene	21.979	228	246616	14.179	ng	99
87) Indeno(1,2,3-cd)pyrene	29.300	276	116788	5.455	ng	# 87
88) Benzo(b)fluoranthene	24.265	252	295914	16.302	ng	# 98
89) Benzo(k)fluoranthene	24.330	252	119356m	6.512	ng	
90) Benzo(a)pyrene	25.193	252	247704	13.853	ng	100
92) Benzo(g,h,i)perylene	30.546	276	113756	6.561	ng	98

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

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