

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055477.D
 Acq On : 5 Nov 2022 19:38
 Operator : CG/JU
 Sample : N5435-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PTS-TP-COMP-1103

Quant Time: Nov 07 00:58:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:51:08 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 11/07/2022
 Supervised By :Jagrut Upadhyay 11/07/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	31313	20.000	ng	0.00
21) Naphthalene-d8	11.056	136	128710	20.000	ng	0.00
39) Acenaphthene-d10	14.845	164	87528	20.000	ng	0.00
64) Phenanthrene-d10	17.577	188	200579	20.000	ng	0.00
76) Chrysene-d12	21.849	240	207308	20.000	ng	0.00
86) Perylene-d12	25.221	264	225625	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.756	112	226873	130.115	ng	0.00
7) Phenol-d6	7.384	99	301544	124.709	ng	0.00
23) Nitrobenzene-d5	9.405	82	187366	77.583	ng	0.00
42) 2,4,6-Tribromophenol	16.326	330	158002	132.341	ng	0.00
45) 2-Fluorobiphenyl	13.470	172	457309	77.465	ng	0.00
79) Terphenyl-d14	20.151	244	764151	71.677	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.619	178	55320	4.903	ng	99
75) Fluoranthene	19.610	202	112182	8.266	ng	99
78) Pyrene	19.969	202	114899	7.857	ng	99
81) Benzo(a)anthracene	21.831	228	69815	4.906	ng	97
83) Chrysene	21.896	228	81378m	5.994	ng	
87) Indeno(1,2,3-cd)pyrene	29.087	276	46323	2.557	ng	# 83
88) Benzo(b)fluoranthene	24.140	252	88484	6.174	ng	97
89) Benzo(k)fluoranthene	24.205	252	35277m	2.449	ng	
90) Benzo(a)pyrene	25.057	252	65372	5.275	ng	98
92) Benzo(g,h,i)perylene	30.304	276	49116	3.304	ng	97

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

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