

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041422\
 Data File : BG053159.D
 Acq On : 14 Apr 2022 16:36
 Operator : CG/JU
 Sample : N2381-01DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 26572DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/15/2022
 Supervised By : Jagrut Upadhyay 04/15/2022

Quant Time: Apr 15 02:21:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.123	152	43494	20.000	ng	0.00
21) Naphthalene-d8	10.931	136	169281	20.000	ng	0.00
39) Acenaphthene-d10	14.743	164	107809	20.000	ng	0.00
64) Phenanthrene-d10	17.487	188	219549	20.000	ng	0.00
76) Chrysene-d12	21.769	240	231649	20.000	ng	0.00
86) Perylene-d12	25.076	264	258333	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.691	112	10287	4.054	ng	0.00
7) Phenol-d6	7.283	99	14997	3.833	ng	0.00
23) Nitrobenzene-d5	9.280	82	11528	2.765	ng	0.00
42) 2,4,6-Tribromophenol	16.230	330	6498	3.788	ng	0.00
45) 2-Fluorobiphenyl	13.369	172	24648	3.153	ng	0.00
79) Terphenyl-d14	20.089	244	40195	2.931	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.537	202	290119	18.873	ng	99
78) Pyrene	19.901	202	384824	23.325	ng	98
81) Benzo(a)anthracene	21.752	228	254799	16.087	ng	98
83) Chrysene	21.822	228	263894m	17.953	ng	
87) Indeno(1,2,3-cd)pyrene	28.842	276	129708	6.758	ng	# 78
88) Benzo(b)fluoranthene	24.031	252	452981	27.791	ng	97
89) Benzo(k)fluoranthene	24.084	252	137622m	8.591	ng	
90) Benzo(a)pyrene	24.918	252	227786	16.405	ng	98
91) Dibenzo(a,h)anthracene	28.877	278	34751m	2.200	ng	
92) Benzo(g,h,i)perylene	30.017	276	104864	6.735	ng	98

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

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