

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051922\
 Data File : BG053630.D
 Acq On : 19 May 2022 17:34
 Operator : CG/JU
 Sample : N2824-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SAC-1

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 05/20/2022
 Supervised By :Jagrut Upadhyay 05/20/2022

Quant Time: May 20 03:57:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 20 03:55:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.076	152	32954	20.000	ng	0.00
21) Naphthalene-d8	10.884	136	127825	20.000	ng	0.00
39) Acenaphthene-d10	14.702	164	87874	20.000	ng	0.00
64) Phenanthrene-d10	17.446	188	169789	20.000	ng	0.00
76) Chrysene-d12	21.758	240	161888	20.000	ng	# 0.02
86) Perylene-d12	25.112	264	151243	20.000	ng	0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.644	112	162418	83.645	ng	0.00
7) Phenol-d6	7.236	99	252638	85.730	ng	0.00
23) Nitrobenzene-d5	9.239	82	184308	61.378	ng	0.00
42) 2,4,6-Tribromophenol	16.189	330	101502	78.317	ng	0.00
45) 2-Fluorobiphenyl	13.322	172	428625	71.277	ng	0.00
79) Terphenyl-d14	20.060	244	545327	62.738	ng	0.00
Target Compounds						Qvalue
83) Chrysene	21.781	228	28258m	2.930	ng	

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

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