

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022822\
 Data File : BG052499.D
 Acq On : 28 Feb 2022 13:35
 Operator : CG/JU
 Sample : N1664-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-6

Quant Time: Mar 01 00:59:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.220	152	29939	20.000	ng	0.00
21) Naphthalene-d8	11.052	136	117273	20.000	ng	#-0.02
39) Acenaphthene-d10	14.864	164	67290	20.000	ng	0.00
64) Phenanthrene-d10	17.613	188	206538	20.000	ng	0.00
76) Chrysene-d12	21.931	240	184184	20.000	ng	0.00
86) Perylene-d12	25.397	264	26283	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.753	112	261380	152.159	ng	0.00
7) Phenol-d6	7.368	99	314124	118.516	ng	0.00
23) Nitrobenzene-d5	9.389	82	261401	96.875	ng	0.00
42) 2,4,6-Tribromophenol	16.350	330	195947	202.691	ng	0.00
45) 2-Fluorobiphenyl	13.489	172	582447	120.275	ng	0.00
79) Terphenyl-d14	20.210	244	1180806	113.210	ng	0.00

Target Compounds	Qvalue

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

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