

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040522\
 Data File : BG053040.D
 Acq On : 5 Apr 2022 20:30
 Operator : CG/JU
 Sample : N2246-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-8-9-3-1

Quant Time: Apr 06 04:15:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 14:24:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	29447	20.000	ng	-0.02
21) Naphthalene-d8	10.936	136	111547	20.000	ng	#-0.02
39) Acenaphthene-d10	14.743	164	81207	20.000	ng	-0.02
64) Phenanthrene-d10	17.486	188	205703	20.000	ng	-0.02
76) Chrysene-d12	21.769	240	250182	20.000	ng	-0.03
86) Perylene-d12	25.064	264	265049	20.000	ng	#-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.691	112	234010	138.395	ng	0.00
7) Phenol-d6	7.277	99	305987	120.033	ng	-0.01
23) Nitrobenzene-d5	9.286	82	262683	118.194	ng	-0.02
42) 2,4,6-Tribromophenol	16.229	330	193637	177.453	ng	-0.02
45) 2-Fluorobiphenyl	13.374	172	551692	100.196	ng	-0.02
79) Terphenyl-d14	20.095	244	1299952	92.429	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.185	163	13800	2.287	ng	Qvalue 99

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

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