

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030822\
 Data File : BG052611.D
 Acq On : 8 Mar 2022 19:18
 Operator : CG/JU
 Sample : N1730-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 31-32-SAMPLE-LOCATION-16

Quant Time: Mar 09 04:19:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 08 01:09:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.212	152	26492	20.000	ng	0.00
21) Naphthalene-d8	11.050	136	104678	20.000	ng	#-0.01
39) Acenaphthene-d10	14.862	164	72731	20.000	ng	0.00
64) Phenanthrene-d10	17.611	188	181864	20.000	ng	0.00
76) Chrysene-d12	21.923	240	204021	20.000	ng	-0.01
86) Perylene-d12	25.383	264	204988	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.745	112	140796	90.546	ng	0.00
7) Phenol-d6	7.361	99	194893	89.198	ng	-0.01
23) Nitrobenzene-d5	9.381	82	137656	59.951	ng	-0.02
42) 2,4,6-Tribromophenol	16.348	330	96560	92.378	ng	0.00
45) 2-Fluorobiphenyl	13.482	172	300578	56.987	ng	0.00
79) Terphenyl-d14	20.202	244	564316	49.532	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.298	163	29610	5.286	ng	Qvalue # 98

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

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