

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060122\
 Data File : BP010658.D
 Acq On : 01 Jun 2022 18:57
 Operator : CG/JU
 Sample : N3076-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-671

Quant Time: Jun 02 10:06:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	114733	20.000	ng	-0.01
21) Naphthalene-d8	10.657	136	492680	20.000	ng	-0.01
39) Acenaphthene-d10	14.492	164	334012	20.000	ng	-0.01
64) Phenanthrene-d10	17.251	188	740940	20.000	ng	0.00
76) Chrysene-d12	21.339	240	814985	20.000	ng	# 0.00
86) Perylene-d12	23.745	264	855840	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	952890	151.514	ng	0.00
7) Phenol-d6	7.040	99	1189137	133.186	ng	0.00
23) Nitrobenzene-d5	9.022	82	1021755	98.049	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	632037	167.167	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	2203420	93.625	ng	-0.01
79) Terphenyl-d14	19.810	244	4255086	98.067	ng	0.00

Target Compounds	Qvalue

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

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