

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060222\
 Data File : BP010673.D
 Acq On : 02 Jun 2022 16:14
 Operator : CG/JU
 Sample : N2990-05RE
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P028-SS001-0006-01RE

Quant Time: Jun 03 00:52:34 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	7.863	152	127593	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	494862	20.000	ng	-0.01
39) Acenaphthene-d10	14.498	164	297131	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	612120	20.000	ng	0.00
76) Chrysene-d12	21.339	240	598209	20.000	ng	0.00
86) Perylene-d12	23.751	264	611790	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	1210214	173.035	ng	0.00
7) Phenol-d6	7.040	99	1575934	158.718	ng	0.00
23) Nitrobenzene-d5	9.022	82	1057643	101.046	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	632448	188.039	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	2159553	103.151	ng	0.00
79) Terphenyl-d14	19.810	244	3448818	108.289	ng	0.00

Target Compounds	Qvalue

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

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