

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005774.D
 Acq On : 08 Jun 2021 11:00
 Operator : CG/JU
 Sample : M2589-39
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(2-4)

Quant Time: Jun 08 11:55:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 07 12:13:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	101008	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	387691	20.000	ng	-0.01
39) Acenaphthene-d10	14.586	164	250680	20.000	ng	-0.02
64) Phenanthrene-d10	17.328	188	579145	20.000	ng	#-0.02
76) Chrysene-d12	21.392	240	758952	20.000	ng	#-0.02
86) Perylene-d12	23.821	264	828856	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	864743	133.102	ng	0.00
7) Phenol-d6	7.134	99	964350	108.883	ng	0.00
23) Nitrobenzene-d5	9.134	82	730781	104.829	ng	0.00
42) 2,4,6-Tribromophenol	16.075	330	624796	194.371	ng	-0.02
45) 2-Fluorobiphenyl	13.222	172	1640918	86.629	ng	-0.01
79) Terphenyl-d14	19.874	244	3204973	79.240	ng	-0.01

Target Compounds	Qvalue

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

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