

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005807.D
 Acq On : 09 Jun 2021 11:34
 Operator : CG/JU
 Sample : M2589-32
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B10(2-4)

Quant Time: Jun 09 12:26:04 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 09 11:54:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	72612	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	270326	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	166603	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	346729	20.000	ng	0.00
76) Chrysene-d12	21.392	240	404259	20.000	ng	0.00
86) Perylene-d12	23.816	264	469371	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.517	112	585622	129.415	ng	0.00
7) Phenol-d6	7.122	99	626387	106.796	ng	0.00
23) Nitrobenzene-d5	9.122	82	471906	85.586	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	380308	133.020	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	1052308	82.941	ng	0.00
79) Terphenyl-d14	19.875	244	1846957	85.549	ng	0.00

Target Compounds	Qvalue

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

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