

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005852.D
 Acq On : 11 Jun 2021 00:58
 Operator : CG/JU
 Sample : PB136995TB
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB136995TB

Quant Time: Jun 11 07:02:34 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	7.946	152	68431	20.000	ng	-0.01
21) Naphthalene-d8	10.751	136	243070	20.000	ng	-0.01
39) Acenaphthene-d10	14.575	164	148343	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	309720	20.000	ng	-0.01
76) Chrysene-d12	21.386	240	352184	20.000	ng	0.00
86) Perylene-d12	23.810	264	436264	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	577617	135.445	ng	0.00
7) Phenol-d6	7.116	99	668791	120.993	ng	0.00
23) Nitrobenzene-d5	9.110	82	411910	83.082	ng	-0.01
42) 2,4,6-Tribromophenol	16.063	330	329893	129.589	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	948109	83.927	ng	-0.01
79) Terphenyl-d14	19.863	244	1570746	83.513	ng	0.00

Target Compounds	Qvalue

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

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