

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062321\
 Data File : BP006041.D
 Acq On : 23 Jun 2021 18:44
 Operator : CG/JU
 Sample : M2789-31
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BASEMENT-WALLS

Quant Time: Jun 24 07:43:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 22 17:52:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	83351	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	319744	20.000	ng	0.00
39) Acenaphthene-d10	14.551	164	198137	20.000	ng	0.00
64) Phenanthrene-d10	17.298	188	404296	20.000	ng	0.00
76) Chrysene-d12	21.369	240	414910	20.000	ng	0.00
86) Perylene-d12	23.780	264	474309	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	255809	49.247	ng	0.01
7) Phenol-d6	7.099	99	503032	74.715	ng	0.00
23) Nitrobenzene-d5	9.087	82	374386	57.406	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	94618	27.827	ng	0.00
45) 2-Fluorobiphenyl	13.181	172	926852	61.426	ng	0.00
79) Terphenyl-d14	19.845	244	1521684	68.674	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.340	178	50970	2.099	ng	99
75) Fluoranthene	19.298	202	59770	2.027	ng	94
84) Bis(2-ethylhexyl)phtha...	21.275	149	710191	40.636	ng	99

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

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