

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061221\
 Data File : BP005898.D
 Acq On : 12 Jun 2021 20:29
 Operator : CG/JU
 Sample : M2677-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 REACTOR

Quant Time: Jun 14 02:10:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 13:53:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	75997	20.000	ng	0.00
21) Naphthalene-d8	10.740	136	276721	20.000	ng	-0.01
39) Acenaphthene-d10	14.563	164	165337	20.000	ng	-0.01
64) Phenanthrene-d10	17.310	188	340772	20.000	ng	-0.01
76) Chrysene-d12	21.381	240	401439	20.000	ng	-0.01
86) Perylene-d12	23.792	264	466258	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	536392	113.256	ng	0.00
7) Phenol-d6	7.117	99	724560	118.032	ng	0.00
23) Nitrobenzene-d5	9.105	82	485447	86.007	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	255353	89.998	ng	-0.01
45) 2-Fluorobiphenyl	13.193	172	1093200	86.824	ng	-0.01
79) Terphenyl-d14	19.857	244	1817108	84.758	ng	-0.01

Target Compounds	Qvalue

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

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