

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005766.D
 Acq On : 08 Jun 2021 06:27
 Operator : CG/JU
 Sample : M2580-33
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B6(0-6)

Quant Time: Jun 08 07:26:10 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	125806	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	476076	20.000	ng	-0.01
39) Acenaphthene-d10	14.592	164	301607	20.000	ng	-0.01
64) Phenanthrene-d10	17.333	188	676333	20.000	ng	-0.01
76) Chrysene-d12	21.398	240	804358	20.000	ng	#-0.02
86) Perylene-d12	23.821	264	832677	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	970469	119.932	ng	0.00
7) Phenol-d6	7.134	99	1070955	97.085	ng	0.00
23) Nitrobenzene-d5	9.134	82	844860	98.694	ng	0.00
42) 2,4,6-Tribromophenol	16.075	330	721311	186.506	ng	-0.02
45) 2-Fluorobiphenyl	13.222	172	1945970	85.387	ng	-0.01
79) Terphenyl-d14	19.875	244	3177056	74.116	ng	-0.01
Target Compounds						
						Qvalue
32) Benzoic acid	10.057	122	22332	9.988	ng	96
50) Dimethylphthalate	14.045	163	125391	5.227	ng	99

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

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