

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005740.D
 Acq On : 07 Jun 2021 13:47
 Operator : CG/JU
 Sample : M2580-18
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B1(0-2)

Quant Time: Jun 07 16:28:45 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	7.975	152	162947	20.000	ng	0.00
21) Naphthalene-d8	10.786	136	597717	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	355650	20.000	ng	0.00
64) Phenanthrene-d10	17.345	188	699711	20.000	ng	# 0.00
76) Chrysene-d12	21.409	240	560034	20.000	ng	# 0.00
86) Perylene-d12	23.845	264	497137	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	1241782	118.482	ng	0.01
7) Phenol-d6	7.145	99	1333208	93.311	ng	0.00
23) Nitrobenzene-d5	9.139	82	1068652	99.431	ng	0.00
42) 2,4,6-Tribromophenol	16.086	330	762808	167.265	ng	0.00
45) 2-Fluorobiphenyl	13.233	172	2262681	84.197	ng	0.00
79) Terphenyl-d14	19.886	244	2968413	99.459	ng	0.00
Target Compounds						
32) Benzoic acid	10.057	122	1906	5.352	ng	# 56
50) Dimethylphthalate	14.051	163	138255	4.887	ng	99

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

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