

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005750.D
 Acq On : 07 Jun 2021 19:31
 Operator : CG/JU
 Sample : M2580-26
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B3(4-6)

Manual Integrations
 APPROVED

mohammad
 6/8/2021 3:37:19 PM

Quant Time: Jun 08 01:15:58 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	133735	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	526535	20.000	ng	0.00
39) Acenaphthene-d10	14.598	164	341746	20.000	ng	0.00
64) Phenanthrene-d10	17.339	188	768268	20.000	ng	# 0.00
76) Chrysene-d12	21.398	240	878631	20.000	ng	#-0.02
86) Perylene-d12	23.833	264	975738	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	1153763	134.130	ng	0.00
7) Phenol-d6	7.140	99	1301509	110.990	ng	0.00
23) Nitrobenzene-d5	9.140	82	1020648	107.803	ng	0.00
42) 2,4,6-Tribromophenol	16.081	330	857957	195.783	ng	-0.01
45) 2-Fluorobiphenyl	13.228	172	2317026	89.728	ng	0.00
79) Terphenyl-d14	19.880	244	3963461	84.645	ng	0.00
Target Compounds						
32) Benzoic acid	10.081	122	4454m	5.910	ng	Qvalue
50) Dimethylphthalate	14.051	163	137780	5.069	ng	98

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

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