

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005770.D
 Acq On : 08 Jun 2021 08:43
 Operator : CG/JU
 Sample : M2603-26
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FAM4-WC7

Manual Integrations
 APPROVED

mohammad
 6/8/2021 3:37:48 PM

Quant Time: Jun 08 09:19:31 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	123110	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	459529	20.000	ng	-0.01
39) Acenaphthene-d10	14.587	164	284059	20.000	ng	-0.02
64) Phenanthrene-d10	17.334	188	610874	20.000	ng	-0.01
76) Chrysene-d12	21.392	240	612449	20.000	ng	#-0.02
86) Perylene-d12	23.821	264	588054	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1044053	131.851	ng	0.00
7) Phenol-d6	7.134	99	1141927	105.785	ng	0.00
23) Nitrobenzene-d5	9.134	82	866319	104.845	ng	0.00
42) 2,4,6-Tribromophenol	16.075	330	652536	179.146	ng	-0.02
45) 2-Fluorobiphenyl	13.222	172	1877470	87.471	ng	-0.01
79) Terphenyl-d14	19.875	244	3067966	93.997	ng	-0.01
Target Compounds						Qvalue
32) Benzoic acid	10.087	122	3673m	5.861	ng	

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

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