

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005761.D
 Acq On : 08 Jun 2021 03:36
 Operator : CG/JU
 Sample : M2603-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FAM4-WC1

Manual Integrations
 APPROVED

mohammad
 6/8/2021 3:37:32 PM

Quant Time: Jun 08 05:44:44 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	127474	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	458409	20.000	ng	-0.01
39) Acenaphthene-d10	14.592	164	274352	20.000	ng	-0.01
64) Phenanthrene-d10	17.333	188	542061	20.000	ng	-0.01
76) Chrysene-d12	21.392	240	444993	20.000	ng	#-0.02
86) Perylene-d12	23.815	264	467000	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1016600	123.989	ng	0.00
7) Phenol-d6	7.140	99	1096761	98.123	ng	0.00
23) Nitrobenzene-d5	9.134	82	833635	101.136	ng	0.00
42) 2,4,6-Tribromophenol	16.075	330	576642	163.912	ng	-0.02
45) 2-Fluorobiphenyl	13.222	172	1780911	85.908	ng	-0.01
79) Terphenyl-d14	19.875	244	2348064	99.013	ng	-0.01
Target Compounds						
32) Benzoic acid	10.140	122	3161m	5.745	ng	Qvalue

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

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