

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063021\
 Data File : BP006124.D
 Acq On : 01 Jul 2021 00:30
 Operator : CG/JU
 Sample : M2896-01 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FS-01

Manual Integrations
 APPROVED

mohammad
 7/1/2021 2:53:01 PM

Quant Time: Jul 01 03:09:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 30 16:39:37 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	120889	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	412728	20.000	ng	0.00
39) Acenaphthene-d10	14.522	164	231997	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	411609	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	509816	20.000	ng	0.00
86) Perylene-d12	23.763	264	613287	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	67302	8.933	ng	0.00
7) Phenol-d6	7.081	99	73957	7.574	ng	0.00
23) Nitrobenzene-d5	9.057	82	48208	5.727	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	30794	7.735	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	100992	5.716	ng	0.00
79) Terphenyl-d14	19.822	244	153985	5.656	ng	0.00
Target Compounds						
						Qvalue
17) 2-Methylphenol	8.357	107	27758	4.176	ng	98
20) 3+4-Methylphenols	8.693	107	53876	6.001	ng	92
27) 2,4-Dimethylphenol	9.881	122	66207m	10.655	ng	
37) 2-Methylnaphthalene	12.351	142	37882	2.228	ng	100

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

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