

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062018\
 Data File : BF106626.D
 Acq On : 20 Jun 2018 18:18
 Operator : JU/SJ
 Sample : J3525-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TPTW-04

Manual Integrations
 APPROVED

Sohil
 6/21/2018 11:27:59 AM

Quant Time: Jun 21 02:00:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	105259	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	450065	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	219186	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	395302	20.00	ng	0.00
75) Chrysene-d12	14.15	240	301702	20.00	ng	0.00
86) Perylene-d12	15.69	264	298558	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	278618	44.82	ng	0.00
7) Phenol-d6	6.60	99	217909	28.07	ng	-0.01
23) Nitrobenzene-d5	7.52	82	678216	83.75	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	250614	98.65	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1117941	86.61	ng	0.00
78) Terphenyl-d14	13.08	244	1185686	95.20	ng	0.00
Target Compounds						
10) Phenol	6.61	94	590039	66.602	ng	99
17) 2-Methylphenol	7.21	107	14664	2.513	ng	98
20) 3+4-Methylphenols	7.37	107	93256	11.857	ng	90
27) 2,4-Dimethylphenol	7.90	122	23787	3.943	ng	99
31) Naphthalene	8.26	128	45626	2.168	ng	98
32) Benzoic acid	8.03	122	243284m	54.550	ng	
49) Dimethylphthalate	9.71	163	43145	2.625	ng	# 87

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

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