

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062018\
 Data File : BF106629.D
 Acq On : 20 Jun 2018 19:39
 Operator : JU/SJ
 Sample : J3529-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TPTW-04

Manual Integrations
 APPROVED

Sohil
 6/21/2018 11:28:01 AM

Quant Time: Jun 21 02:19:06 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	6.96	152	106018	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	434298	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	221243	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	447281	20.00	ng	0.00
75) Chrysene-d12	14.15	240	298284	20.00	ng	0.00
86) Perylene-d12	15.69	264	300112	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	391988	62.61	ng	0.00
7) Phenol-d6	6.61	99	313899	40.15	ng	0.00
23) Nitrobenzene-d5	7.52	82	669490	85.67	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	369243	144.00	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1138127	87.51	ng	0.00
78) Terphenyl-d14	13.09	244	1277937	103.78	ng	0.00
Target Compounds						
10) Phenol	6.62	94	596700	66.872	ng	99
17) 2-Methylphenol	7.22	107	14072	2.395	ng	93
20) 3+4-Methylphenols	7.37	107	95610	12.097	ng	93
27) 2,4-Dimethylphenol	7.90	122	23607	4.055	ng	96
31) Naphthalene	8.27	128	46917	2.311	ng	100
32) Benzoic acid	8.08	122	244103m	56.493	ng	
49) Dimethylphthalate	9.71	163	111154	6.699	ng	# 94

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

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