Quantitation Report (LSC Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061918\
Data File : BF106593.D
Acq On : 19 Jun 2018 20:16
Operator : JU/SJ
Sample : J3522-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TPTW-01

Quant Time: Jun 20 02:25:05 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	116849	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	439348	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	184982	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	280974	20.00	ng	0.00
75) Chrysene-d12	14.15	240	272814	20.00	ng	0.00
86) Perylene-d12	15.68	264	320356	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	259691	37.63	ng	0.00
7) Phenol-d6	6.60	99	199388	23.14	ng	0.00
23) Nitrobenzene-d5	7.52	82	657344	83.15	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	161352	75.26	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1050741	98.49	ng	0.00
78) Terphenyl-d14	13.08	244	815983	72.45	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.61	94	28741	2.922	ng	95
49) Dimethylphthalate	9.71	163	83828	6.042	ng	99
51) Acenaphthene	10.04	154	47528	4.120	ng	99
57) Fluorene	10.55	166	42938	3.425	ng	98
70) Phenanthrene	11.52	178	73695	5.438	ng	98
72) Carbazole	11.73	167	61220	4.727	ng	99

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

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