

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061118\
 Data File : BF106335.D
 Acq On : 12 Jun 2018 10:19
 Operator : JU/SJ
 Sample : J3398-08
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-31

Quant Time: Jun 12 11:48:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	212500	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	766945	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	329828	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	524520	20.00	ng	0.00
75) Chrysene-d12	14.15	240	508398	20.00	ng	0.00
86) Perylene-d12	15.69	264	344681	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	655124	52.18	ng	0.00
7) Phenol-d6	6.59	99	520410	32.81	ng	0.00
23) Nitrobenzene-d5	7.54	82	1107312	84.94	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	400032	126.05	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1711216	95.86	ng	0.00
78) Terphenyl-d14	13.09	244	1390825	67.95	ng	0.00
Target Compounds						
10) Phenol	6.60	94	75676	4.370	ng	96
17) 2-Methylphenol	7.21	107	48650	4.039	ng	# 87
31) Naphthalene	8.28	128	78533	2.265	ng	# 95
32) Benzoic acid	8.07	122	218786	30.627	ng	90
37) 2-Methylnaphthalene	8.97	142	131108	5.552	ng	97
49) Dimethylphthalate	9.72	163	86056	3.652	ng	# 97

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

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