

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082718\
 Data File : BM016632.D
 Acq On : 28 Aug 2018 08:22
 Operator : SJ/JU
 Sample : J4652-04DL 10X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SONO-WDL

Quant Time: Aug 28 09:27:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 27 13:01:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	8487	20.00	ng	0.03
21) Naphthalene-d8	10.83	136	27862	20.00	ng	0.02
38) Acenaphthene-d10	14.65	164	23290	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	85250	20.00	ng	0.01
75) Chrysene-d12	21.54	240	113336	20.00	ng	0.00
86) Perylene-d12	23.83	264	121610	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	10239	23.64	ng	0.04
7) Phenol-d6	7.22	99	9060	15.81	ng	0.04
23) Nitrobenzene-d5	9.21	82	39996	65.54	ng	0.02
41) 2,4,6-Tribromophenol	16.14	330	68297	104.90	ng	0.00
44) 2-Fluorobiphenyl	13.26	172	151050	62.34	ng	0.00
78) Terphenyl-d14	19.97	244	493033	92.07	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.13	163	6005	2.684	ng	# 78
58) 2,3,4,6-Tetrachlorophenol	15.28	232	253603	343.415	ng	# 100
69) Pentachlorophenol	17.05	266	9014098	11152.564	ng	99

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

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