

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080818\
 Data File : BM016270.D
 Acq On : 09 Aug 2018 05:15
 Operator : SJ/JU
 Sample : J4328-04
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-04

Quant Time: Aug 09 07:08:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 16:23:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	43332	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	152963	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	96130	20.00	ng	0.00
63) Phenanthrene-d10	17.50	188	157958	20.00	ng	0.00
75) Chrysene-d12	21.63	240	168813	20.00	ng	0.00
86) Perylene-d12	23.97	264	182102	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	130437	50.00	ng	0.00
7) Phenol-d6	7.30	99	94307	27.68	ng	0.00
23) Nitrobenzene-d5	9.32	82	299048	90.93	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	128432	105.34	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	553518	70.06	ng	0.00
78) Terphenyl-d14	20.09	244	677602	84.02	ng	0.00
Target Compounds						
8) 2-Chlorophenol	7.69	128	16963	5.471	ng	84
13) 1,4-Dichlorobenzene	8.17	146	108799	29.561	ng	# 95
14) 1,2-Dichlorobenzene	8.49	146	30704	8.526	ng	# 93
20) 3+4-Methylphenols	8.92	107	19364	5.780	ng	# 82
27) 2,4-Dimethylphenol	10.13	122	33853	17.584	ng	# 75
31) Naphthalene	10.99	128	36449	4.708	ng	# 92

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080818\
Data File : BM016270.D
Acq On : 09 Aug 2018 05:15
Operator : SJ/JU
Sample : J4328-04
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MW-04

Quant Time: Aug 09 07:08:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 08 16:23:37 2018
Response via : Initial Calibration

