

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082718\
 Data File : BM016634.D
 Acq On : 28 Aug 2018 09:33
 Operator : SJ/JU
 Sample : J4652-04DL 10X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SONO-WDL

Manual Integrations
 APPROVED

Sohil
 8/28/2018 11:17:02 AM

Quant Time: Aug 28 10:58:20 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	8.00	152	71172	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	261333	20.00	ng	0.00
38) Acenaphthene-d10	14.64	164	207546	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	800111	20.00	ng	0.00
75) Chrysene-d12	21.53	240	1168241	20.00	ng	0.00
86) Perylene-d12	23.81	264	1217990	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	13435m	3.70	ng	0.05
7) Phenol-d6	7.23	99	10538m	2.19	ng	0.05
23) Nitrobenzene-d5	9.23	82	42356m	7.40	ng	0.04
41) 2,4,6-Tribromophenol	16.14	330	56794	9.79	ng	0.00
44) 2-Fluorobiphenyl	13.26	172	127057	5.88	ng	0.00
78) Terphenyl-d14	19.97	244	416368	7.54	ng	0.00
Target Compounds						
58) 2,3,4,6-Tetrachlorophenol	15.28	232	208012	31.609	ng	Qvalue # 100
69) Pentachlorophenol	17.05	266	7218736	951.607	ng	98

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

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