

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020718\
 Data File : BM014172.D
 Acq On : 07 Feb 2018 21:55
 Operator : SJ/JU
 Sample : J1161-03
 Misc : MDL 1 PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-SOIL-03-QT1-2018

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:50:33 PM

Quant Time: Feb 08 11:29:21 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 17:16:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	82833	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	343843	20.00	ng	0.00
38) Acenaphthene-d10	14.20	164	222643	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	538348	20.00	ng	0.00
75) Chrysene-d12	21.17	240	571931	20.00	ng	0.00
86) Perylene-d12	23.34	264	511407	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	622877	131.07	ng	0.00
7) Phenol-d6	6.73	99	834068	123.92	ng	0.00
23) Nitrobenzene-d5	8.70	82	707430	91.90	ng	0.00
41) 2,4,6-Tribromophenol	15.70	330	407309	128.93	ng	0.00
44) 2-Fluorobiphenyl	12.81	172	1594594	89.41	ng	0.00
78) Terphenyl-d14	19.60	244	2546605	87.00	ng	0.00
Target Compounds						
15) Benzyl Alcohol	7.86	79	17953m	2.88	ng	Qvalue
40) Hexachlorocyclopentadiene	12.33	237	653	6.30	ng	# 28
64) 4,6-Dinitro-2-methylphenol	15.42	198	525	5.69	ng	# 28
69) Pentachlorophenol	16.66	266	1228	6.18	ng	# 65

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

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