

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF085981.D
 Acq On : 2 Apr 2016 00:39
 Operator : UM/SJ
 Sample : H1824-03
 Misc : LOD 5PPM
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-SOIL2016-01

Manual Integrations
 APPROVED

Sohil
 4/4/2016 7:39:26 PM

Quant Time: Apr 02 01:40:19 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	102098	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	409635	20.00	ng	-0.01
38) Acenaphthene-d10	9.81	164	190751	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	378102	20.00	ng	-0.01
75) Chrysene-d12	13.93	240	346569	20.00	ng	0.00
86) Perylene-d12	15.33	264	289867	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	766843	128.38	ng	0.00
7) Phenol-d6	6.42	99	1019626	134.39	ng	0.00
23) Nitrobenzene-d5	7.34	82	617713	88.93	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	279215	155.95	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1105745	96.38	ng	0.00
78) Terphenyl-d14	12.88	244	1063393	72.13	ng	0.00
Target Compounds						Qvalue
3) Pyridine	3.29	79	17751m	2.80	ng	
4) n-Nitrosodimethylamine	3.13	42	7879m	2.83	ng	
32) Benzoic acid	7.82	122	7689	1.60	ng	96
40) Hexachlorocyclopentadiene	8.92	237	1765	6.54	ng	90
53) 2,4-Dinitrophenol	9.92	184	1168	3.04	ng	# 70

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

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