

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102418\
 Data File : BN003276.D
 Acq On : 24 Oct 2018 17:55
 Operator : MJ/SJ
 Sample : J5164-01
 Misc : LOD 2PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LOD-MDL-SOIL-01-QT4-2018

Quant Time: Oct 25 06:38:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 24 14:21:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	94555	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	440986	20.00	ng	0.00
39) Acenaphthene-d10	14.24	164	278103	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	661679	20.00	ng	0.00
76) Chrysene-d12	21.21	240	754482	20.00	ng	0.00
87) Perylene-d12	23.42	264	793114	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	767572	141.39	ng	0.00
7) Phenol-d6	6.80	99	1066551	138.16	ng	0.00
23) Nitrobenzene-d5	8.76	82	640807	83.34	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	538018	143.67	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1697341	82.79	ng	0.00
79) Terphenyl-d14	19.64	244	2714061	76.93	ng	0.00
Target Compounds						
25) Isophorone	9.32	82	29197	2.165	ng	99
28) bis(2-Chloroethoxy)methane	9.82	93	17928	1.973	ng	92
30) 1,2,4-Trichlorobenzene	10.25	180	14548	1.931	ng	99
61) 4-Chlorophenyl-phenylether	15.30	204	23242	1.996	ng	97

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

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