

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091919\
 Data File : BP000326.D
 Acq On : 19 Sep 2019 13:41
 Operator : HP/JU
 Sample : K4938-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MG-NW-SOIL

Manual Integrations
 APPROVED

mohammad
 9/20/2019 1:08:35 PM

Quant Time: Sep 20 04:54:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	227337	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	843196	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	518922	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	1041824	20.00	ng	0.00
76) Chrysene-d12	13.99	240	1000255	20.00	ng	0.00
87) Perylene-d12	15.47	264	1059303	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1242953	94.39	ng	0.00
7) Phenol-d6	6.43	99	1617107	100.18	ng	0.00
23) Nitrobenzene-d5	7.35	82	885080	67.72	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	593444	115.30	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2029486	70.39	ng	0.00
79) Terphenyl-d14	12.93	244	2837432	59.51	ng	0.00
Target Compounds						
10) Phenol	6.43	94	50638	2.761	ng	81
50) Dimethylphthalate	9.55	163	319486	9.037	ng	98
75) Fluoranthene	12.55	202	235809	4.206	ng	95
78) Pyrene	12.78	202	213594	2.854	ng	98
88) Benzo(b)fluoranthene	15.04	252	149678m	2.369	ng	

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

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