

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100430.D
 Acq On : 11 Nov 2017 10:33
 Operator : SJ/JU
 Sample : I6368-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-RCA-SOIL

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:54:04 PM

Quant Time: Nov 13 01:37:33 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	180186	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	712183	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	315768	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	525260	20.00	ng	0.00
75) Chrysene-d12	14.06	240	382215	20.00	ng	0.00
86) Perylene-d12	15.56	264	343649	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	320143	28.48	ng	0.00
7) Phenol-d6	6.53	99	1060814	76.53	ng	0.00
23) Nitrobenzene-d5	7.46	82	755139	70.10	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	42453	13.19	ng	-0.01
44) 2-Fluorobiphenyl	9.26	172	1224961	59.07	ng	0.00
78) Terphenyl-d14	13.01	244	891452	53.59	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.53	94	66107	4.543	ng	86
31) Naphthalene	8.20	128	115298	3.361	ng	98
37) 2-Methylnaphthalene	8.89	142	123408	5.419	ng	99
49) Dimethylphthalate	9.65	163	210570	8.734	ng	99
51) Acenaphthene	9.97	154	53670	2.688	ng	97
57) Fluorene	10.49	166	66021	3.220	ng	99
70) Phenanthrene	11.45	178	416756	15.069	ng	98
71) Anthracene	11.50	178	96263	3.431	ng	99
74) Fluoranthene	12.64	202	366821	12.515	ng	96
77) Pyrene	12.87	202	344424	12.641	ng	100
79) Butylbenzylphthalate	13.48	149	58975	5.308	ng	96
80) Benzo(a)anthracene	14.05	228	164283	7.138	ng	95
82) Chrysene	14.09	228	160477	7.200	ng	97
83) Bis(2-ethylhexyl)phthalate	14.04	149	34318	2.348	ng	# 98
85) Indeno(1,2,3-cd)pyrene	17.05	276	51221	2.841	ng	96
87) Benzo(b)fluoranthene	15.12	252	164835m	8.096	ng	
88) Benzo(k)fluoranthene	15.14	252	52602m	2.734	ng	
89) Benzo(a)pyrene	15.49	252	110767	6.137	ng	# 91
91) Benzo(a,h,i)perylene	17.51	276	50019	3.462	ng	# 91

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

