

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011419\
 Data File : BF112053.D
 Acq On : 14 Jan 2019 16:25
 Operator : JU/SJ
 Sample : K1080-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONCRETE-PILE

Manual Integrations
 APPROVED

Sohil
 1/15/2019 12:19:19 PM

Quant Time: Jan 14 21:16:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 11 21:37:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	153088	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	547599	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	292072	20.00	ng	-0.01
64) Phenanthrene-d10	11.39	188	459377	20.00	ng	0.00
76) Chrysene-d12	14.03	240	328556	20.00	ng	-0.01
87) Perylene-d12	15.51	264	339687	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	810374	90.56	ng	0.00
7) Phenol-d6	6.52	99	1158636	100.46	ng	0.00
23) Nitrobenzene-d5	7.43	82	596005	58.17	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	267638	70.45	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	1203928	60.11	ng	0.00
79) Terphenyl-d14	12.97	244	1319288	76.22	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.62	163	114530	5.220	ng	97
71) Phenanthrene	11.42	178	76972	3.134	ng	99
75) Fluoranthene	12.60	202	108243	4.293	ng	99
78) Pyrene	12.83	202	108359	4.793	ng	98
81) Benzo(a)anthracene	14.02	228	43105	2.280	ng	97
83) Chrysene	14.06	228	40387	2.240	ng	96
86) Indeno(1,2,3-cd)pyrene	16.99	276	29758	2.264	ng	94
88) Benzo(b)fluoranthene	15.07	252	60763m	2.990	ng	
90) Benzo(a)pyrene	15.44	252	46453	2.600	ng	96
92) Benzo(g,h,i)perylene	17.44	276	29808	2.121	ng	97

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

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