

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113852.D
 Acq On : 19 Apr 2019 12:18
 Operator : JU/SJ
 Sample : K2453-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-STONES

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:26:28 PM

Quant Time: Apr 19 23:56:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	208497	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	743030	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	342631	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	537973	20.00	ng	0.00
76) Chrysene-d12	14.06	240	412999	20.00	ng	0.02
87) Perylene-d12	15.54	264	561051	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1402776	111.87	ng	0.01
7) Phenol-d6	6.52	99	1795886	114.85	ng	0.00
23) Nitrobenzene-d5	7.46	82	1087154	89.10	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	440881	116.47	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1892505	84.23	ng	0.00
79) Terphenyl-d14	13.01	244	1448619	73.55	ng	0.01
Target Compounds						
50) Dimethylphthalate	9.64	163	262898	10.964	ng	99
71) Phenanthrene	11.44	178	91209	3.329	ng	# 94
75) Fluoranthene	12.63	202	151473	5.216	ng	94
78) Pyrene	12.86	202	146996	5.050	ng	95
81) Benzo(a)anthracene	14.06	228	87492	3.624	ng	94
83) Chrysene	14.09	228	90869	3.792	ng	96
85) Di-n-octyl phthalate	14.70	149	281957	14.088	ng	# 67
86) Indeno(1,2,3-cd)pyrene	17.02	276	72266	3.694	ng	95
88) Benzo(b)fluoranthene	15.11	252	135697m	3.880	ng	
90) Benzo(a)pyrene	15.48	252	99211	3.208	ng	# 40
92) Benzo(g,h,i)perylene	17.47	276	67441	2.411	ng	# 91

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

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