

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040919\
 Data File : BF113667.D
 Acq On : 9 Apr 2019 18:01
 Operator : JU/SJ
 Sample : K2338-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-DEEP

Manual Integrations
 APPROVED

Sohil
 4/10/2019 6:20:39 PM

Quant Time: Apr 10 06:48:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	113532	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	432209	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	212381	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	428155	20.00	ng	-0.01
76) Chrysene-d12	13.82	240	317780	20.00	ng	-0.02
87) Perylene-d12	15.23	264	333371	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	830344	124.72	ng	0.00
7) Phenol-d6	6.33	99	1091332	132.99	ng	0.00
23) Nitrobenzene-d5	7.25	82	649258	87.99	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	342935	134.42	ng	-0.01
45) 2-Fluorobiphenyl	9.03	172	1154228	87.49	ng	-0.02
79) Terphenyl-d14	12.77	244	1177027	75.41	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.43	163	179420	11.461	ng	99
75) Fluoranthene	12.40	202	84447	3.704	ng	96
78) Pyrene	12.63	202	79049	3.269	ng	97
81) Benzo(a)anthracene	13.82	228	41566	2.267	ng	99
83) Chrysene	13.85	228	47051	2.532	ng	99
86) Indeno(1,2,3-cd)pyrene	16.60	276	36922	2.152	ng	96
88) Benzo(b)fluoranthene	14.84	252	66459m	3.257	ng	
90) Benzo(a)pyrene	15.17	252	46721	2.577	ng	98

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

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