

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021918\
 Data File : BF103083.D
 Acq On : 19 Feb 2018 15:38
 Operator : SJ/JU
 Sample : J1604-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4

Manual Integrations
 APPROVED

Sohil
 2/20/2018 10:20:30 AM

Quant Time: Feb 19 17:40:35 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	166457	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	699751	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	295287	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	479203	20.00	ng	0.00
75) Chrysene-d12	14.05	240	304145	20.00	ng	0.00
86) Perylene-d12	15.54	264	320818	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1073121	97.91	ng	0.01
7) Phenol-d6	6.51	99	1290075	97.75	ng	0.00
23) Nitrobenzene-d5	7.44	82	833119	82.59	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	249202	104.85	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1247100	70.08	ng	0.00
78) Terphenyl-d14	12.99	244	914074	71.16	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	82868	3.599	ng	99
70) Phenanthrene	11.43	178	234558	8.789	ng	99
74) Fluoranthene	12.62	202	254375	9.473	ng	97
77) Pyrene	12.85	202	203119	8.517	ng	97
80) Benzo(a)anthracene	14.04	228	68906	3.602	ng	98
82) Chrysene	14.07	228	78266	4.059	ng	100
85) Indeno(1,2,3-cd)pyrene	17.04	276	38765	2.424	ng	95
87) Benzo(b)fluoranthene	15.10	252	98429m	4.732	ng	
89) Benzo(a)pyrene	15.47	252	67335	3.596	ng	# 95
91) Benzo(g,h,i)perylene	17.50	276	35656	2.397	ng	94

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

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