

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103704.D
 Acq On : 16 Mar 2018 18:53
 Operator : JU/SJ
 Sample : J1757-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-031418-B

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:01:03 PM

Quant Time: Mar 17 02:07:17 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	178369	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	720290	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	319079	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	497581	20.00	ng	0.00
75) Chrysene-d12	14.15	240	325808	20.00	ng	0.00
86) Perylene-d12	15.68	264	315112	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1136576	96.92	ng	0.01
7) Phenol-d6	6.59	99	1422113	99.12	ng	0.00
23) Nitrobenzene-d5	7.53	82	858247	83.09	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	279407	101.84	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1274207	63.70	ng	0.00
78) Terphenyl-d14	13.09	244	945616	67.37	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	94089	3.865	ng	98
70) Phenanthrene	11.52	178	124691	4.581	ng	99
74) Fluoranthene	12.72	202	106724	3.806	ng	99
77) Pyrene	12.95	202	122971	5.139	ng	99
80) Benzo(a)anthracene	14.14	228	50184m	2.433	ng	
82) Chrysene	14.17	228	50060	2.546	ng	96
87) Benzo(b)fluoranthene	15.22	252	49847m	2.504	ng	
89) Benzo(a)pyrene	15.61	252	40740	2.256	ng	# 80

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

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