

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100217\
 Data File : BF099016.D
 Acq On : 2 Oct 2017 22:52
 Operator : SJ/JU
 Sample : I5532-11 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-66-1-2

Manual Integrations
 APPROVED

Sohil
 10/3/2017 6:51:03 PM

Quant Time: Oct 03 04:19:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	128943	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	498892	20.00	ng	-0.02
38) Acenaphthene-d10	9.95	164	190264	20.00	ng	-0.02
63) Phenanthrene-d10	11.43	188	292772	20.00	ng	-0.01
75) Chrysene-d12	14.07	240	263314	20.00	ng	-0.01
86) Perylene-d12	15.57	264	228749	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	274729	33.92	ng	-0.02
7) Phenol-d6	6.53	99	533302	53.37	ng	-0.02
23) Nitrobenzene-d5	7.46	82	323306	41.56	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	7071	4.54	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	581066	50.98	ng	-0.02
78) Terphenyl-d14	13.02	244	419716	36.25	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.65	163	93548	6.514	ng	99
70) Phenanthrene	11.46	178	130776	8.345	ng	96
74) Fluoranthene	12.65	202	138435	8.724	ng	99
77) Pyrene	12.87	202	124395	6.195	ng	98
80) Benzo(a)anthracene	14.06	228	56069	3.517	ng	97
82) Chrysene	14.10	228	58686	3.866	ng	96
87) Benzo(b)fluoranthene	15.12	252	66921m	4.758	ng	
89) Benzo(a)pyrene	15.50	252	41397	3.207	ng	# 82

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

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