

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099942.D
 Acq On : 29 Oct 2017 4:40
 Operator : SJ/JU
 Sample : I6059-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 101817-TIDO-F-D-B

Quant Time: Oct 30 05:17:04 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 15:58:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	107568	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	418605	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	172507	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	244275	20.00	ng	-0.02
75) Chrysene-d12	13.96	240	182719	20.00	ng	-0.03
86) Perylene-d12	15.42	264	176595	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	279073	40.73	ng	0.00
7) Phenol-d6	6.43	99	237924	29.29	ng	-0.01
23) Nitrobenzene-d5	7.36	82	568047	87.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	98737	62.81	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	1058940	94.71	ng	-0.01
78) Terphenyl-d14	12.90	244	688979	82.40	ng	-0.02
Target Compounds						
10) Phenol	6.45	94	24522	2.749	ng	93
49) Dimethylphthalate	9.54	163	70322	5.148	ng	99
51) Acenaphthene	9.86	154	30597	2.869	ng	96
70) Phenanthrene	11.34	178	70812	5.137	ng	96

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

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