

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099946.D
 Acq On : 29 Oct 2017 6:32
 Operator : SJ/JU
 Sample : I6100-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 102317-TIDO-F-D-B

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:20:47 PM

Quant Time: Oct 30 05:30:44 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	112585	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	441100	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	174464	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	244527	20.00	ng	-0.02
75) Chrysene-d12	13.96	240	193348	20.00	ng	-0.03
86) Perylene-d12	15.42	264	184155	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	294832	41.11	ng	0.00
7) Phenol-d6	6.43	99	219544	25.82	ng	-0.01
23) Nitrobenzene-d5	7.36	82	572095	83.50	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	100667	63.32	ng	-0.02
44) 2-Fluorobiphenyl	9.14	172	1080215	95.53	ng	-0.01
78) Terphenyl-d14	12.90	244	695612	78.62	ng	-0.02
Target Compounds						
10) Phenol	6.45	94	19625	2.102	ng	90
49) Dimethylphthalate	9.54	163	31412	2.274	ng	97
70) Phenanthrene	11.34	178	32545	2.358	ng	95

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

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