

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
 Data File : BF100872.D
 Acq On : 23 Nov 2017 13:45
 Operator : SJ/JU
 Sample : I6567-04
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112017-TIDO-F-U-S

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:21:12 PM

Quant Time: Nov 25 01:00:28 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 15:22:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	73236	20.00	ng	-0.02
21) Naphthalene-d8	8.14	136	287295	20.00	ng	-0.03
38) Acenaphthene-d10	9.89	164	130116	20.00	ng	-0.03
63) Phenanthrene-d10	11.38	188	227724	20.00	ng	-0.02
75) Chrysene-d12	14.02	240	141308	20.00	ng	-0.02
86) Perylene-d12	15.49	264	156333	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	184980	40.49	ng	-0.02
7) Phenol-d6	6.48	99	137321	24.37	ng	-0.03
23) Nitrobenzene-d5	7.42	82	454507	104.59	ng	-0.03
41) 2,4,6-Tribromophenol	10.68	330	122712	92.55	ng	-0.03
44) 2-Fluorobiphenyl	9.22	172	877673	102.71	ng	-0.02
78) Terphenyl-d14	12.97	244	614282	99.88	ng	-0.02
Target Compounds						
10) Phenol	6.49	94	15335	2.593	ng	95
49) Dimethylphthalate	9.60	163	24807	2.497	ng	99

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

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