

Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
 Data File : BF097979.D
 Acq On : 23 Aug 2017 3:14
 Operator : SJ/JU
 Sample : I4914-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 289-COMP

Manual Integrations
 APPROVED

Sohil
 8/23/2017 3:04:02 PM

Quant Time: Aug 23 05:11:32 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	126791	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	484938	20.00	ng	0.00
38) Acenaphthene-d10	9.78	164	195286	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	309679	20.00	ng	0.00
75) Chrysene-d12	13.90	240	267850	20.00	ng	0.00
86) Perylene-d12	15.32	264	218636	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	932186	111.83	ng	0.01
7) Phenol-d6	6.39	99	1256543	110.95	ng	0.00
23) Nitrobenzene-d5	7.32	82	782394	87.65	ng	0.00
41) 2,4,6-Tribromophenol	10.57	330	152790	90.17	ng	0.00
44) 2-Fluorobiphenyl	9.11	172	1038032	78.09	ng	0.00
78) Terphenyl-d14	12.85	244	682133	53.11	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.50	163	164211	11.504	ng	98
74) Fluoranthene	12.47	202	59949	3.481	ng	98
77) Pyrene	12.70	202	53154	2.426	ng	96
83) Bis(2-ethylhexyl)phthalate	13.89	149	131886	14.170	ng	99
87) Benzo(b)fluoranthene	14.91	252	29923m	2.171	ng	

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

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