

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100109.D
 Acq On : 3 Nov 2017 3:12
 Operator : SJ/JU
 Sample : I6274-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110117-TIDO-F-D-S

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:20:50 AM

Quant Time: Nov 03 04:44:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	80847	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	331318	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	147623	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	239532	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	144638	20.00	ng	-0.02
86) Perylene-d12	15.41	264	131911	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	213748	41.51	ng	0.00
7) Phenol-d6	6.42	99	178932	29.30	ng	-0.01
23) Nitrobenzene-d5	7.34	82	425930	82.77	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	102228	75.99	ng	-0.02
44) 2-Fluorobiphenyl	9.13	172	836243	87.40	ng	0.00
78) Terphenyl-d14	12.88	244	622811	94.10	ng	-0.02
Target Compounds						
31) Naphthalene	8.07	128	61632	3.854	ng	99
70) Phenanthrene	11.32	178	35880	2.654	ng	98

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

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