

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101917\
 Data File : BF099688.D
 Acq On : 20 Oct 2017 6:38
 Operator : SJ/JU
 Sample : I5927-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 101317-TIDO-E-D-M

Quant Time: Oct 20 14:02:11 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 20 13:11:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	91554	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	372624	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	161978	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	239069	20.00	ng	0.00
75) Chrysene-d12	13.98	240	156908	20.00	ng	0.00
86) Perylene-d12	15.44	264	169124	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	269447	46.85	ng	0.00
7) Phenol-d6	6.46	99	202001	28.47	ng	0.00
23) Nitrobenzene-d5	7.39	82	564315	97.12	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	119896	90.46	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1071244	110.41	ng	0.00
78) Terphenyl-d14	12.92	244	680299	98.60	ng	0.00
Target Compounds						
51) Acenaphthene	9.89	154	27733	2.736	ng	98
74) Fluoranthene	12.56	202	45090	3.480	ng	96
77) Pyrene	12.79	202	28551	2.386	ng	98

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

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