

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100116.D
 Acq On : 3 Nov 2017 6:21
 Operator : SJ/JU
 Sample : I6273-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 103117-TIDO-E-D-B

Quant Time: Nov 03 07:38:03 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	89450	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	364697	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	166060	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	284252	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	179719	20.00	ng	-0.02
86) Perylene-d12	15.41	264	151191	20.00	ng	-0.05

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.40	112	286098	50.21	ng	0.00
7) Phenol-d6	6.42	99	242320	35.87	ng	-0.01
23) Nitrobenzene-d5	7.35	82	541758	95.64	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	142290	94.02	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	1018945	94.67	ng	0.00
78) Terphenyl-d14	12.88	244	893870	108.69	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.07	128	373858	21.237	ng	99
37) 2-Methylnaphthalene	8.76	142	61197	5.187	ng	97
51) Acenaphthene	9.84	154	70108	6.828	ng	97
54) Dibenzofuran	10.02	168	70165	4.841	ng	95
57) Fluorene	10.36	166	59309	4.943	ng	99
70) Phenanthrene	11.32	178	178543	11.130	ng	97
74) Fluoranthene	12.52	202	55882	3.352	ng	98
77) Pyrene	12.75	202	35703	2.613	ng	99

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

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