

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\
 Data File : BF097457.D
 Acq On : 8 Aug 2017 5:04
 Operator : SJ/JU
 Sample : I4545-01
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-09-BASE

Quant Time: Aug 08 07:45:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.41	152	96481	20.00	ng	0.00
21) Naphthalene-d8	7.69	136	356010	20.00	ng	0.00
38) Acenaphthene-d10	9.43	164	138154	20.00	ng	0.00
63) Phenanthrene-d10	10.91	188	224911	20.00	ng	0.00
75) Chrysene-d12	13.53	240	211675	20.00	ng	0.00
86) Perylene-d12	14.87	264	142391	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.00	112	377636	59.00	ng	0.00
7) Phenol-d6	6.08	99	458106	62.70	ng	0.00
23) Nitrobenzene-d5	6.98	82	300178	49.46	ng	-0.01
41) 2,4,6-Tribromophenol	10.23	330	63164	57.87	ng	0.00
44) 2-Fluorobiphenyl	8.77	172	358653	44.06	ng	0.00
78) Terphenyl-d14	12.49	244	238913	23.01	ng	0.00
Target Compounds						
11) bis(2-Chloroethyl)ether	6.16	93	19417	2.833	ng	88
13) 1,4-Dichlorobenzene	6.43	146	18497	2.531	ng	89
14) 1,2-Dichlorobenzene	6.58	146	15352	2.406	ng	94
16) 2,2'-oxybis(1-Chloropropan	6.71	45	28690	2.087	ng	95
27) 2,4-Dimethylphenol	7.40	122	11995	2.127	ng	99
31) Naphthalene	7.72	128	206581	12.310	ng	99
37) 2-Methylnaphthalene	8.40	142	87661	8.250	ng	97
49) Dimethylphthalate	9.17	163	75742	7.220	ng	98
51) Acenaphthene	9.47	154	20225	2.576	ng	98
70) Phenanthrene	10.93	178	42388	3.517	ng	98

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

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