

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097059.D
 Acq On : 26 Jul 2017 6:06
 Operator : SJ/JU
 Sample : I4399-02 5X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-N7-BASE

Quant Time: Jul 26 06:56:45 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	113470	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	428893	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	171363	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	249704	20.00	ng	0.00
75) Chrysene-d12	13.62	240	201280	20.00	ng	0.00
86) Perylene-d12	14.97	264	145721	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.10	112	141092	18.74	ng	0.00
7) Phenol-d6	6.16	99	168627	19.62	ng	-0.01
23) Nitrobenzene-d5	7.06	82	97677	13.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.31	330	25276	18.67	ng	-0.01
44) 2-Fluorobiphenyl	8.86	172	155020	15.35	ng	-0.01
78) Terphenyl-d14	12.58	244	78679	7.97	ng	0.00
Target Compounds						
11) bis(2-Chloroethyl)ether	6.23	93	69230	8.588	ng	86
12) 1,3-Dichlorobenzene	6.44	146	58117	6.700	ng	97
13) 1,4-Dichlorobenzene	6.52	146	181218	21.082	ng	94
14) 1,2-Dichlorobenzene	6.67	146	241834	32.222	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.79	45	128343	7.938	ng	88
17) 2-Methylphenol	6.78	107	17218	2.772	ng	# 78
20) 3+4-Methylphenols	6.94	107	18045	2.224	ng	# 58
27) 2,4-Dimethylphenol	7.46	122	166739	24.537	ng	97
31) Naphthalene	7.82	128	3021105	149.430	ng	99
37) 2-Methylnaphthalene	8.50	142	635170	49.620	ng	99
49) Dimethylphthalate	9.26	163	34254	2.632	ng	99

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

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