

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080417\
 Data File : BF097389.D
 Acq On : 5 Aug 2017 7:22
 Operator : SJ/JU
 Sample : I4588-02 10X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-M8-ESW

Manual Integrations
 APPROVED

Sohil
 8/7/2017 6:10:24 PM

Quant Time: Aug 07 12:22:27 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 01:17:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.44	152	104557	20.00	ng	-0.06
21) Naphthalene-d8	7.72	136	383251	20.00	ng	-0.06
38) Acenaphthene-d10	9.46	164	153205	20.00	ng	-0.07
63) Phenanthrene-d10	10.93	188	208090	20.00	ng	-0.07
75) Chrysene-d12	13.56	240	169073	20.00	ng	-0.07
86) Perylene-d12	14.90	264	131655	20.00	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	49949	7.20	ng	-0.04
7) Phenol-d6	6.13	99	66464	8.39	ng	-0.04
23) Nitrobenzene-d5	7.01	82	38278	5.86	ng	-0.06
41) 2,4,6-Tribromophenol	10.25	330	8899	7.35	ng	-0.07
44) 2-Fluorobiphenyl	8.80	172	69124	7.66	ng	-0.07
78) Terphenyl-d14	12.52	244	38619	4.66	ng	-0.07
Target Compounds						
11) bis(2-Chloroethyl)ether	6.18	93	29650	3.992	ng	Qvalue 94
13) 1,4-Dichlorobenzene	6.46	146	24552	3.100	ng	91
14) 1,2-Dichlorobenzene	6.61	146	49603	7.172	ng	95
20) 3+4-Methylphenols	6.91	107	84989	11.369	ng	# 67
22) Acetophenone	6.86	105	45189	4.910	ng	# 98
27) 2,4-Dimethylphenol	7.42	122	37246	6.134	ng	97
37) 2-Methylnaphthalene	8.43	142	23897	2.089	ng	# 94
39) 1,2,4,5-Tetrachlorobenzene	8.59	216	15579	3.811	ng	# 76
48) Acenaphthylene	9.33	152	34917	2.362	ng	# 90
77) Pyrene	12.36	202	28963	2.092	ng	# 94
87) Benzo(b)fluoranthene	14.55	252	34307m	4.335	ng	
91) Benzo(a,h,i)perylene	16.46	276	12954	2.080	ng	# 90

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

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