

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101221\
 Data File : BN016900.D
 Acq On : 12 Oct 2021 16:20
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4EC

Quant Time: Oct 12 16:55:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 12 02:31:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	19805	0.400	ng	-0.02
7) Naphthalene-d8	10.445	136	87157	0.400	ng	-0.01
13) Acenaphthene-d10	14.294	164	46213	0.400	ng	0.00
19) Phenanthrene-d10	17.030	188	85966	0.400	ng	#-0.01
29) Chrysene-d12	21.238	240	83289	0.400	ng	0.00
35) Perylene-d12	23.481	264	82920	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.282	112	17606	0.347	ng	-0.02
5) Phenol-d6	6.859	99	19156	0.337	ng	0.00
8) Nitrobenzene-d5	8.810	82	19986	0.339	ng	-0.03
11) 2-Methylnaphthalene-d10	12.034	152	48461	0.384	ng	0.00
14) 2,4,6-Tribromophenol	15.779	330	6906	0.308	ng	-0.01
15) 2-Fluorobiphenyl	12.911	172	68811	0.389	ng	-0.01
27) Fluoranthene-d10	19.068	212	82291	0.365	ng	0.00
31) Terphenyl-d14	19.679	244	72785	0.384	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.263	88	5910	0.399	ng	# 92
3) n-Nitrosodimethylamine	3.595	42	6269	0.387	ng	98
6) bis(2-Chloroethyl)ether	7.117	93	21760	0.391	ng	100
9) Naphthalene	10.495	128	89757	0.378	ng	99
10) Hexachlorobutadiene	10.774	225	17880	0.401	ng	# 99
12) 2-Methylnaphthalene	12.105	142	57796	0.389	ng	99
16) Acenaphthylene	14.002	152	69859	0.354	ng	100
17) Acenaphthene	14.358	154	50340	0.384	ng	100
18) Fluorene	15.335	166	60534	0.383	ng	100
20)						

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