

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092220\
 Data File : BN011873.D
 Acq On : 23 Sep 2020 04:54
 Operator : CG/JU
 Sample : L4082-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE105D1-20200915

Quant Time: Sep 23 08:30:04 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 14:36:55 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	3764	0.40	ng	0.00
7) Naphthalene-d8	10.478	136	15802	0.40	ng	0.00
13) Acenaphthene-d10	14.331	164	8486	0.40	ng	0.00
19) Phenanthrene-d10	17.078	188	16844	0.40	ng	0.00
29) Chrysene-d12	21.258	240	15313	0.40	ng	#-0.01
36) Perylene-d12	23.484	264	14969	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.299	112	1305	0.12	ng	0.00
5) Phenol-d6	6.869	99	961	0.07	ng	0.00
8) Nitrobenzene-d5	8.843	82	3628	0.23	ng	0.00
11) 2-Methylnaphthalene-d10	12.064	152	5009	0.18	ng	0.00
14) 2,4,6-Tribromophenol	15.816	330	1464	0.34	ng	0.00
15) 2-Fluorobiphenyl	12.948	172	9927	0.27	ng	0.00
27) Fluoranthene-d10	19.107	212	12048	0.23	ng	0.00
31) Terphenyl-d14	19.707	244	12633	0.36	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.261	88	17061	3.05	ng	Qvalue 97
6) bis(2-Chloroethyl)ether	7.127	93	357	0.03	ng	96
9) Naphthalene	10.516	128	1039	0.02	ng	# 78
34) Bis(2-ethylhexyl)phtha...	21.184	149	3459	0.13	ng	92

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

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