

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121620\
 Data File : BN013157.D
 Acq On : 16 Dec 2020 07:43
 Operator : CG/JU
 Sample : L4996-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE132D7-20201207

Quant Time: Dec 16 08:43:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 15 16:49:20 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.346	152	718	0.40	ng	0.00
7) Naphthalene-d8	10.125	136	2323	0.40	ng	# 0.01
13) Acenaphthene-d10	14.004	164	1476	0.40	ng	-0.01
19) Phenanthrene-d10	16.788	188	3159	0.40	ng	# 0.01
29) Chrysene-d12	21.016	240	3447	0.40	ng	# 0.01
36) Perylene-d12	23.099	264	3917	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.970	112	353	0.12	ng	0.00
5) Phenol-d6	6.549	99	236	0.08	ng	0.00
8) Nitrobenzene-d5	8.528	82	610	0.25	ng	0.01
11) 2-Methylnaphthalene-d10	11.733	152	1062	0.26	ng	0.00
14) 2,4,6-Tribromophenol	15.539	330	217	0.23	ng	0.00
15) 2-Fluorobiphenyl	12.633	172	1580	0.26	ng	0.00
27) Fluoranthene-d10	18.826	212	3875	0.37	ng	0.00
31) Terphenyl-d14	19.442	244	3461	0.40	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.933	88	2059	1.64	ng	# 93
6) bis(2-Chloroethyl)ether	6.798	93	66	0.04	ng	94

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

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