

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN113022\
 Data File : BN022983.D
 Acq On : 30 Nov 2022 13:18
 Operator : CG/JU
 Sample : N5671-13DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP04-20221115DL

Quant Time: Nov 30 23:00:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN112522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 30 00:45:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.064	152	6344	0.400	ng	0.00
7) Naphthalene-d8	10.883	136	16517	0.400	ng	0.00
13) Acenaphthene-d10	14.698	164	7461	0.400	ng	0.00
19) Phenanthrene-d10	17.439	188	14428	0.400	ng	0.00
29) Chrysene-d12	21.625	240	9222	0.400	ng	# 0.00
35) Perylene-d12	24.118	264	7179	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.594	112	665	0.040	ng	0.00
5) Phenol-d6	7.204	99	535	0.025	ng	0.00
8) Nitrobenzene-d5	9.217	82	1825	0.142	ng	-0.01
11) 2-Methylnaphthalene-d10	12.461	152	3235	0.090	ng	0.00
14) 2,4,6-Tribromophenol	16.186	330	12	0.003	ng	0.00
15) 2-Fluorobiphenyl	13.330	172	5146	0.151	ng	0.00
27) Fluoranthene-d10	19.473	212	4900	0.125	ng	0.00
31) Terphenyl-d14	20.067	244	3843	0.177	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.442	88	19046	2.365	ng	Qvalue 99

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

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