

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102622\
 Data File : BN022334.D
 Acq On : 26 Oct 2022 13:30
 Operator : CG/JU
 Sample : N5231-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP-100-20221020

Quant Time: Oct 27 05:58:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 27 05:55:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	3803	0.400	ng	0.00
7) Naphthalene-d8	10.773	136	12877	0.400	ng	0.00
13) Acenaphthene-d10	14.621	164	7380	0.400	ng	0.01
19) Phenanthrene-d10	17.375	188	13114	0.400	ng	# 0.00
29) Chrysene-d12	21.582	240	6989	0.400	ng	0.00
35) Perylene-d12	24.043	264	5816	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.479	112	1230	0.140	ng	0.00
5) Phenol-d6	7.119	99	843	0.074	ng	0.01
8) Nitrobenzene-d5	9.129	82	2881	0.278	ng	0.01
11) 2-Methylnaphthalene-d10	12.372	152	6082	0.257	ng	0.00
14) 2,4,6-Tribromophenol	16.121	330	583	0.207	ng	0.00
15) 2-Fluorobiphenyl	13.242	172	9777	0.303	ng	0.00
27) Fluoranthene-d10	19.411	212	9938	0.285	ng	0.00
31) Terphenyl-d14	20.006	244	7925	0.564	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.284	88	17238	3.503	ng	Qvalue 99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102622\
Data File : BN022334.D
Acq On : 26 Oct 2022 13:30
Operator : CG/JU
Sample : N5231-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SP-100-20221020

Quant Time: Oct 27 05:58:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 27 05:55:40 2022
Response via : Initial Calibration

