

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092922\
 Data File : BN021958.D
 Acq On : 29 Sep 2022 17:11
 Operator : CG/JU
 Sample : N4846-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP09-20220923

Quant Time: Sep 30 02:45:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 30 02:27:15 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.033	152	4097	0.400	ng	0.00
7) Naphthalene-d8	10.855	136	12407	0.400	ng	0.00
13) Acenaphthene-d10	14.686	164	6303	0.400	ng	0.00
19) Phenanthrene-d10	17.437	188	12100	0.400	ng	0.00
29) Chrysene-d12	21.636	240	8600	0.400	ng	0.00
35) Perylene-d12	24.134	264	6230	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.563	112	811	0.074	ng	0.00
5) Phenol-d6	7.188	99	617	0.045	ng	0.00
8) Nitrobenzene-d5	9.211	82	1792	0.173	ng	0.00
11) 2-Methylnaphthalene-d10	12.448	152	5234	0.196	ng	0.00
14) 2,4,6-Tribromophenol	16.184	330	573	0.212	ng	0.00
15) 2-Fluorobiphenyl	13.317	172	5611	0.201	ng	0.00
27) Fluoranthene-d10	19.471	212	8564	0.261	ng	0.00
31) Terphenyl-d14	20.060	244	7469	0.393	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.382	88	3134	0.540	ng	# 92
34) Bis(2-ethylhexyl)phtha...	21.529	149	3059	0.153	ng	# 99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092922\
Data File : BN021958.D
Acq On : 29 Sep 2022 17:11
Operator : CG/JU
Sample : N4846-03
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP09-20220923

Quant Time: Sep 30 02:45:43 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 30 02:27:15 2022
Response via : Initial Calibration

