

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102922\
 Data File : BM037273.D
 Acq On : 29 Oct 2022 16:27
 Operator : CG/JU
 Sample : N5219-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BACKYARD-SURFACE-A

Quant Time: Oct 31 02:06:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 31 01:34:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	457771	20.000	ng	0.00
21) Naphthalene-d8	10.674	136	1792376	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	1006406	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1858236	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1370477	20.000	ng	0.00
86) Perylene-d12	23.779	264	1359300	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	2690988	101.559	ng	0.00
7) Phenol-d6	7.045	99	3492578	97.119	ng	0.00
23) Nitrobenzene-d5	9.045	82	2095958	58.998	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	975247	82.229	ng	0.00
45) 2-Fluorobiphenyl	13.133	172	4290160	56.640	ng	0.00
79) Terphenyl-d14	19.868	244	4797434	62.775	ng	0.00

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102922\
Data File : BM037273.D
Acq On : 29 Oct 2022 16:27
Operator : CG/JU
Sample : N5219-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BACKYARD-SURFACE-A

Quant Time: Oct 31 02:06:29 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 31 01:34:15 2022
Response via : Initial Calibration

