

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
 Data File : BM030337.D
 Acq On : 09 Jun 2021 11:47
 Operator : CG/JU
 Sample : M2632-01RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PRODUCT01-IW-PS-D1BRRE

Quant Time: Jun 09 13:40:37 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	206611	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	527271	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	35007	20.000	ng	# 0.00
64) Phenanthrene-d10	17.204	188	166007	20.000	ng	# 0.00
76) Chrysene-d12	21.362	240	620	20.000	ng	# 0.00
86) Perylene-d12	23.633	264	762	20.000	ng	#-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1287924	101.337	ng	0.00
7) Phenol-d6	7.022	99	64621	3.512	ng	0.01
23) Nitrobenzene-d5	9.004	82	1563749	144.484	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	1088332	2472.724	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3649579	1370.988	ng	0.00
79) Terphenyl-d14	19.815	244	256361	7262.624	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.352	88	17036	3.101	ng	# 76
46) 1,1'-Biphenyl	13.310	154	8659	3.054	ng	87

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
Data File : BM030337.D
Acq On : 09 Jun 2021 11:47
Operator : CG/JU
Sample : M2632-01RE
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PRODUCT01-IW-PS-D1BRRE

Quant Time: Jun 09 13:40:37 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 08 20:51:08 2021
Response via : Initial Calibration

