Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
Data File : BN034143.D
Acq On : 22 Sep 2024 02:59
Operator : JU/RC
Sample : P4111-01
Misc :
ALS Vial : 108 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RW7-SP100-20240918

Quant Time: Sep 23 00:02:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 16:09:28 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.423	152	5853	0.400	ng	-0.01
7) Naphthalene-d8	10.180	136	15081	0.400	ng	#-0.02
13) Acenaphthene-d10	14.072	164	8155	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	15584	0.400	ng	0.00
29) Chrysene-d12	21.044	240	10906	0.400	ng	#-0.01
35) Perylene-d12	23.168	264	10163	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	2238	0.135	ng	0.00
5) Phenol-d6	6.629	99	1277	0.066	ng	0.00
8) Nitrobenzene-d5	8.568	82	3496	0.303	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	7034	0.316	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	806	0.218	ng	0.00
15) 2-Fluorobiphenyl	12.683	172	11756	0.354	ng	-0.02
27) Fluoranthene-d10	18.876	212	14317	0.364	ng	-0.01
31) Terphenyl-d14	19.494	244	10670	0.490	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.068	88	48318	6.603	ng	# 87
6) bis(2-Chloroethyl)ether	6.874	93	585	0.034	ng	92

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

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