Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
Data File : BN034484.D
Acq On : 09 Oct 2024 01:20
Operator : RC/JU
Sample : P4226-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP07-20240924

Quant Time: Oct 09 01:44:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 26 14:44:02 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.380	152	4467	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	12487	0.400	ng	-0.01
13) Acenaphthene-d10	14.026	164	7076	0.400	ng	-0.02
19) Phenanthrene-d10	16.790	188	12271	0.400	ng	-0.01
29) Chrysene-d12	21.006	240	8090	0.400	ng	0.00
35) Perylene-d12	23.116	264	9497	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.025	112	1563	0.123	ng	0.00
5) Phenol-d6	6.585	99	751	0.051	ng	-0.01
8) Nitrobenzene-d5	8.536	82	2041	0.214	ng	0.00
11) 2-Methylnaphthalene-d10	11.734	152	4690	0.254	ng	-0.02
14) 2,4,6-Tribromophenol	15.549	330	852	0.266	ng	0.00
15) 2-Fluorobiphenyl	12.647	172	6847	0.238	ng	-0.01
27) Fluoranthene-d10	18.832	212	11306	0.365	ng	-0.01
31) Terphenyl-d14	19.449	244	7938	0.491	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	24613	4.407	ng	100
6) bis(2-Chloroethyl)ether	6.824	93	450	0.034	ng	91

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

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