

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022324\
 Data File : BN029956.D
 Acq On : 23 Feb 2024 14:52
 Operator : MA/JU
 Sample : PB159108BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB159108BL

Quant Time: Feb 23 17:10:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN012924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 23:52:05 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.862	152	7409	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	19550	0.400	ng	#-0.02
13) Acenaphthene-d10	14.500	164	9100	0.400	ng	-0.01
19) Phenanthrene-d10	17.255	188	17655	0.400	ng	0.00
29) Chrysene-d12	21.456	240	13204	0.400	ng	# 0.00
35) Perylene-d12	23.820	264	13757	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.435	112	7935	0.446	ng	0.00
5) Phenol-d6	7.031	99	8758	0.426	ng	0.00
8) Nitrobenzene-d5	9.025	82	5652	0.347	ng	-0.01
11) 2-Methylnaphthalene-d10	12.246	152	8582	0.322	ng	-0.01
14) 2,4,6-Tribromophenol	16.002	330	1020	0.349	ng	0.00
15) 2-Fluorobiphenyl	13.121	172	12500	0.320	ng	-0.01
27) Fluoranthene-d10	19.289	212	13353	0.312	ng	0.00
31) Terphenyl-d14	19.898	244	10301	0.314	ng	0.00

Target Compounds Qvalue

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

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