

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102423\
 Data File : BN028372.D
 Acq On : 24 Oct 2023 11:21
 Operator : MA/JU
 Sample : PB156426BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB156426BL

Quant Time: Oct 25 01:03:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN101223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 15:14:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	4694	0.400	ng	0.00
7) Naphthalene-d8	10.693	136	15825	0.400	ng	-0.01
13) Acenaphthene-d10	14.523	164	8472	0.400	ng	-0.01
19) Phenanthrene-d10	17.281	188	15198	0.400	ng	#-0.01
29) Chrysene-d12	21.480	240	11784	0.400	ng	# 0.00
35) Perylene-d12	23.928	264	12659	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.423	112	6250	0.429	ng	0.00
5) Phenol-d6	7.055	99	7082	0.395	ng	0.00
8) Nitrobenzene-d5	9.091	82	5331	0.380	ng	-0.01
11) 2-Methylnaphthalene-d10	12.279	152	9901	0.416	ng	-0.01
14) 2,4,6-Tribromophenol	16.040	330	1410	0.341	ng	-0.02
15) 2-Fluorobiphenyl	13.144	172	13174	0.442	ng	-0.02
27) Fluoranthene-d10	19.319	212	15270	0.376	ng	0.00
31) Terphenyl-d14	19.909	244	12336	0.493	ng	0.00

Target Compounds Qvalue

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

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