

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034468.D
 Acq On : 08 Oct 2024 13:08
 Operator : RC/JU
 Sample : PB163839BL
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB163839BL

Quant Time: Oct 08 14:09:47 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	5371	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	14280	0.400	ng	-0.01
13) Acenaphthene-d10	14.040	164	7333	0.400	ng	0.00
19) Phenanthrene-d10	16.808	188	12040	0.400	ng	0.00
29) Chrysene-d12	21.017	240	10785	0.400	ng	0.00
35) Perylene-d12	23.130	264	11785	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.025	112	5382	0.353	ng	0.00
5) Phenol-d6	6.593	99	3908	0.220	ng	0.00
8) Nitrobenzene-d5	8.536	82	2920	0.267	ng	0.00
11) 2-Methylnaphthalene-d10	11.746	152	6810	0.323	ng	0.00
14) 2,4,6-Tribromophenol	15.567	330	881	0.265	ng	0.02
15) 2-Fluorobiphenyl	12.661	172	9232	0.310	ng	0.00
27) Fluoranthene-d10	18.839	212	11316	0.373	ng	0.00
31) Terphenyl-d14	19.461	244	7393	0.343	ng	0.00

Target Compounds Qvalue

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

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