

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011924\
 Data File : BN029297.D
 Acq On : 19 Jan 2024 10:28
 Operator : MA/JU
 Sample : PB158469BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB158469BL

Quant Time: Jan 19 11:03:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN011824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 18 16:49:42 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.941	152	4202	0.400	ng	0.00
7) Naphthalene-d8	10.744	136	11478	0.400	ng	0.00
13) Acenaphthene-d10	14.575	164	6193	0.400	ng	0.00
19) Phenanthrene-d10	17.330	188	15223	0.400	ng	0.00
29) Chrysene-d12	21.528	240	12580	0.400	ng	# 0.00
35) Perylene-d12	23.938	264	12949	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.514	112	3878	0.372	ng	0.00
5) Phenol-d6	7.103	99	5007	0.362	ng	0.00
8) Nitrobenzene-d5	9.099	82	3043	0.286	ng	0.00
11) 2-Methylnaphthalene-d10	12.329	152	5462	0.279	ng	0.00
14) 2,4,6-Tribromophenol	16.064	330	603	0.317	ng	0.00
15) 2-Fluorobiphenyl	13.207	172	8810	0.306	ng	0.00
27) Fluoranthene-d10	19.359	212	12640	0.274	ng	0.00
31) Terphenyl-d14	19.968	244	10293	0.326	ng	0.00

Target Compounds Qvalue

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

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