

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110624\
 Data File : BN034881.D
 Acq On : 06 Nov 2024 19:52
 Operator : RC/JU
 Sample : PB164229BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB164229BL

Quant Time: Nov 07 00:43:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN103024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 05 00:03:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	5724	0.400	ng	-0.01
7) Naphthalene-d8	10.340	136	15408	0.400	ng	-0.01
13) Acenaphthene-d10	14.208	164	6105	0.400	ng	-0.01
19) Phenanthrene-d10	16.952	188	10732	0.400	ng	#-0.02
29) Chrysene-d12	21.149	240	4888	0.400	ng	#-0.01
35) Perylene-d12	23.315	264	4283	0.400	ng	#-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.199	112	5644	0.326	ng	-0.01
5) Phenol-d6	6.759	99	7243	0.321	ng	0.00
8) Nitrobenzene-d5	8.717	82	4762	0.393	ng	-0.01
11) 2-Methylnaphthalene-d10	11.939	152	7627	0.342	ng	-0.01
14) 2,4,6-Tribromophenol	15.698	330	346	0.115	ng	-0.02
15) 2-Fluorobiphenyl	12.829	172	10842	0.451	ng	-0.01
27) Fluoranthene-d10	18.990	212	8736	0.347	ng	-0.01
31) Terphenyl-d14	19.603	244	5007	0.477	ng	-0.01

Target Compounds Qvalue

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

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