

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110824\
 Data File : BN034904.D
 Acq On : 08 Nov 2024 12:24
 Operator : RC/JU
 Sample : PB164235BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB164235BL

Quant Time: Nov 08 12:49:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 15:02:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	5987	0.400	ng	0.00
7) Naphthalene-d8	10.340	136	17079	0.400	ng	0.00
13) Acenaphthene-d10	14.208	164	7385	0.400	ng	0.00
19) Phenanthrene-d10	16.952	188	14844	0.400	ng	# 0.00
29) Chrysene-d12	21.149	240	8034	0.400	ng	0.00
35) Perylene-d12	23.315	264	6658	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.199	112	6158	0.369	ng	0.00
5) Phenol-d6	6.759	99	8119	0.367	ng	0.00
8) Nitrobenzene-d5	8.707	82	5118	0.384	ng	0.00
11) 2-Methylnaphthalene-d10	11.935	152	8613	0.370	ng	0.00
14) 2,4,6-Tribromophenol	15.698	330	498	0.268	ng	0.00
15) 2-Fluorobiphenyl	12.829	172	12718	0.408	ng	0.00
27) Fluoranthene-d10	18.990	212	13244	0.396	ng	0.00
31) Terphenyl-d14	19.598	244	7850	0.522	ng	0.00

Target Compounds Qvalue

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

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