

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102622\
 Data File : BN022335.D
 Acq On : 26 Oct 2022 14:09
 Operator : CG/JU
 Sample : N5231-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP-201-20221020

Quant Time: Oct 27 05:59:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 27 05:55:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	4038	0.400	ng	0.00
7) Naphthalene-d8	10.773	136	13082	0.400	ng	0.00
13) Acenaphthene-d10	14.621	164	6679	0.400	ng	0.01
19) Phenanthrene-d10	17.375	188	11112	0.400	ng	# 0.00
29) Chrysene-d12	21.582	240	7839	0.400	ng	0.00
35) Perylene-d12	24.043	264	6865	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.479	112	1417	0.151	ng	0.00
5) Phenol-d6	7.119	99	986	0.081	ng	0.01
8) Nitrobenzene-d5	9.129	82	3005	0.286	ng	0.01
11) 2-Methylnaphthalene-d10	12.372	152	5768	0.240	ng	0.00
14) 2,4,6-Tribromophenol	16.121	330	587	0.230	ng	0.00
15) 2-Fluorobiphenyl	13.242	172	8424	0.289	ng	0.00
27) Fluoranthene-d10	19.411	212	9998	0.338	ng	0.00
31) Terphenyl-d14	20.006	244	7958	0.505	ng	0.00

Target Compounds Qvalue

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

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