

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081021\
 Data File : BN015837.D
 Acq On : 10 Aug 2021 13:25
 Operator : CG/JU
 Sample : PB137774BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB138188BL

Quant Time: Aug 11 10:53:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN080921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 09 19:13:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	9820	0.400	ng	0.00
7) Naphthalene-d8	10.723	136	48320	0.400	ng	0.00
13) Acenaphthene-d10	14.560	164	27912	0.400	ng	0.00
19) Phenanthrene-d10	17.297	188	56573	0.400	ng	#-0.01
29) Chrysene-d12	21.493	240	60057	0.400	ng	# 0.00
35) Perylene-d12	23.922	264	56256	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.485	112	9363	0.366	ng	0.00
5) Phenol-d6	7.080	99	12424	0.359	ng	0.00
8) Nitrobenzene-d5	9.076	82	14038	0.370	ng	0.00
11) 2-Methylnaphthalene-d10	12.314	152	33283	0.429	ng	0.00
14) 2,4,6-Tribromophenol	16.044	330	3499	0.280	ng	0.00
15) 2-Fluorobiphenyl	13.190	172	44332	0.395	ng	0.00
27) Fluoranthene-d10	19.336	212	76789	0.447	ng	0.00
31) Terphenyl-d14	19.936	244	80227	0.455	ng	0.00

Target Compounds Qvalue

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

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