

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090921\
 Data File : BN016290.D
 Acq On : 10 Sep 2021 14:51
 Operator : CG/JU
 Sample : PB138988BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB138988BL

Quant Time: Sep 10 15:38:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 10 14:34:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	12202	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	61377	0.400	ng	0.00
13) Acenaphthene-d10	14.497	164	33493	0.400	ng	0.01
19) Phenanthrene-d10	17.237	188	67775	0.400	ng	0.00
29) Chrysene-d12	21.429	240	52042	0.400	ng	# 0.00
35) Perylene-d12	23.820	264	28038	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.475	112	11241	0.361	ng	0.00
5) Phenol-d6	7.052	99	12484	0.320	ng	0.00
8) Nitrobenzene-d5	9.025	82	14859	0.301	ng	0.00
11) 2-Methylnaphthalene-d10	12.247	152	33375	0.348	ng	0.00
14) 2,4,6-Tribromophenol	15.994	330	3408	0.248	ng	0.01
15) 2-Fluorobiphenyl	13.114	172	42622	0.325	ng	0.00
27) Fluoranthene-d10	19.271	212	69071	0.347	ng	0.00
31) Terphenyl-d14	19.867	244	55977	0.429	ng	0.00

Target Compounds	Qvalue

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

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