

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090421\
 Data File : BN016207.D
 Acq On : 05 Sep 2021 00:44
 Operator : CG/JU
 Sample : PB138734BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB138734BL

Quant Time: Sep 06 06:20:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 06 02:08:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.873	152	14522	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	77002	0.400	ng	0.00
13) Acenaphthene-d10	14.497	164	44070	0.400	ng	0.00
19) Phenanthrene-d10	17.237	188	83069	0.400	ng	#-0.01
29) Chrysene-d12	21.429	240	84565	0.400	ng	# 0.00
35) Perylene-d12	23.816	264	80776	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.475	112	14159	0.345	ng	0.02
5) Phenol-d6	7.043	99	16808	0.321	ng	0.00
8) Nitrobenzene-d5	9.025	82	21512	0.348	ng	0.00
11) 2-Methylnaphthalene-d10	12.252	152	43739	0.357	ng	0.00
14) 2,4,6-Tribromophenol	15.981	330	5564	0.283	ng	-0.01
15) 2-Fluorobiphenyl	13.127	172	60615	0.365	ng	0.00
27) Fluoranthene-d10	19.276	212	86412	0.346	ng	0.00
31) Terphenyl-d14	19.872	244	82740	0.387	ng	0.00

Target Compounds Qvalue

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

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