

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092021\
 Data File : BN016338.D
 Acq On : 20 Sep 2021 10:44
 Operator : CG/JU
 Sample : PB139183BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139183BL

Quant Time: Sep 20 11:14:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 20 11:09:23 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.688	152	24666	0.400	ng	0.00
7) Naphthalene-d8	10.457	136	112262	0.400	ng	0.00
13) Acenaphthene-d10	14.307	164	60066	0.400	ng	0.00
19) Phenanthrene-d10	17.054	188	118156	0.400	ng	0.00
29) Chrysene-d12	21.259	240	116169	0.400	ng	# 0.00
35) Perylene-d12	23.522	264	96562	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.328	112	24163	0.390	ng	0.02
5) Phenol-d6	6.877	99	27216	0.363	ng	0.00
8) Nitrobenzene-d5	8.822	82	26432	0.396	ng	0.00
11) 2-Methylnaphthalene-d10	12.047	152	71227	0.403	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	5324	0.302	ng	0.00
15) 2-Fluorobiphenyl	12.937	172	97539	0.415	ng	0.00
27) Fluoranthene-d10	19.098	212	139448	0.408	ng	0.00
31) Terphenyl-d14	19.708	244	122886	0.457	ng	0.00

Target Compounds Qvalue

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

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