

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091721\
 Data File : BN016308.D
 Acq On : 17 Sep 2021 21:40
 Operator : CG/JU
 Sample : PB139166BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139166BL

Quant Time: Sep 18 00:18:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 17 23:59:14 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.688	152	25641	0.400	ng	0.00
7) Naphthalene-d8	10.457	136	120249	0.400	ng	# 0.00
13) Acenaphthene-d10	14.307	164	67849	0.400	ng	0.00
19) Phenanthrene-d10	17.066	188	127497	0.400	ng	# 0.00
29) Chrysene-d12	21.259	240	119170	0.400	ng	# 0.00
35) Perylene-d12	23.526	264	112870	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.328	112	22299	0.347	ng	0.03
5) Phenol-d6	6.877	99	26051	0.334	ng	0.00
8) Nitrobenzene-d5	8.835	82	22748	0.318	ng	0.00
11) 2-Methylnaphthalene-d10	12.052	152	66866	0.353	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	4923	0.247	ng	0.00
15) 2-Fluorobiphenyl	12.937	172	90198	0.340	ng	0.00
27) Fluoranthene-d10	19.103	212	125417	0.340	ng	0.00
31) Terphenyl-d14	19.713	244	111736	0.405	ng	0.00

Target Compounds Qvalue

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

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