

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120221\
 Data File : BN017683.D
 Acq On : 02 Dec 2021 15:05
 Operator : CG/JU
 Sample : PB141111BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB141111BL

Quant Time: Dec 02 15:41:04 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN113021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 02 01:51:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.727	152	18184	0.400	ng	0.00
7) Naphthalene-d8	10.510	136	69870	0.400	ng	# 0.00
13) Acenaphthene-d10	14.359	164	46154	0.400	ng	0.00
19) Phenanthrene-d10	17.092	188	92650	0.400	ng	#-0.01
29) Chrysene-d12	21.293	240	93170	0.400	ng	0.01
35) Perylene-d12	23.592	264	95417	0.400	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.348	112	13222	0.320	ng	0.00
5) Phenol-d6	6.906	99	11882	0.272	ng	0.00
8) Nitrobenzene-d5	8.862	82	11947	0.403	ng	-0.01
11) 2-Methylnaphthalene-d10	12.102	152	34746	0.285	ng	0.00
14) 2,4,6-Tribromophenol	15.843	330	4061	0.265	ng	0.00
15) 2-Fluorobiphenyl	12.976	172	56114	0.312	ng	-0.01
27) Fluoranthene-d10	19.134	212	80374	0.277	ng	0.00
31) Terphenyl-d14	19.740	244	69254	0.339	ng	0.00

Target Compounds Qvalue

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

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