

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111721\
 Data File : BN017490.D
 Acq On : 17 Nov 2021 10:14
 Operator : CG/JU
 Sample : PB140437BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB140437BL

Quant Time: Nov 17 10:59:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN111621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 16 14:06:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.874	152	25817	0.400	ng	0.00
7) Naphthalene-d8	10.661	136	113776	0.400	ng	# 0.00
13) Acenaphthene-d10	14.498	164	57873	0.400	ng	0.00
19) Phenanthrene-d10	17.238	188	100007	0.400	ng	0.00
29) Chrysene-d12	21.420	240	68635	0.400	ng	# 0.00
35) Perylene-d12	23.794	264	49124	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.477	112	19814	0.322	ng	0.03
5) Phenol-d6	7.044	99	20142	0.305	ng	0.00
8) Nitrobenzene-d5	9.014	82	20482	0.290	ng	0.00
11) 2-Methylnaphthalene-d10	12.253	152	54788	0.293	ng	0.00
14) 2,4,6-Tribromophenol	15.983	330	4048	0.275	ng	0.00
15) 2-Fluorobiphenyl	13.128	172	71206	0.326	ng	0.00
27) Fluoranthene-d10	19.267	212	83508	0.280	ng	0.00
31) Terphenyl-d14	19.868	244	63469	0.377	ng	0.00

Target Compounds Qvalue

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

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