

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021221\
 Data File : BN013613.D
 Acq On : 12 Feb 2021 14:38
 Operator : CG/JU
 Sample : PB134610BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB134610BL

Quant Time: Feb 12 17:18:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN020821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 08 15:13:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.555	152	6003	0.40	ng	0.00
7) Naphthalene-d8	10.315	136	23099	0.40	ng	# 0.00
13) Acenaphthene-d10	14.181	164	11884	0.40	ng	0.00
19) Phenanthrene-d10	16.934	188	23715	0.40	ng	0.00
29) Chrysene-d12	21.133	240	19105	0.40	ng	0.00
36) Perylene-d12	23.304	264	17767	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.188	112	4857	0.37	ng	0.00
5) Phenol-d6	6.734	99	5469	0.34	ng	0.00
8) Nitrobenzene-d5	8.693	82	4384	0.34	ng	0.00
11) 2-Methylnaphthalene-d10	11.911	152	11696	0.32	ng	0.00
14) 2,4,6-Tribromophenol	15.691	330	617	0.18	ng	0.01
15) 2-Fluorobiphenyl	12.798	172	16000	0.36	ng	0.00
27) Fluoranthene-d10	18.975	212	18884	0.30	ng	0.00
31) Terphenyl-d14	19.586	244	16976	0.40	ng	0.00

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

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