

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121021\
 Data File : BN017871.D
 Acq On : 10 Dec 2021 15:13
 Operator : CG/JU
 Sample : PB141233BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB141233BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/11/2021
 Supervised By : Yogesh Patel 12/15/2021

Quant Time: Dec 11 01:47:48 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN120721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 08 01:18:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.715	152	19645	0.400	ng	0.00
7) Naphthalene-d8	10.507	136	70102	0.400	ng	0.01
13) Acenaphthene-d10	14.347	164	44286	0.400	ng	0.00
19) Phenanthrene-d10	17.117	188	77255	0.400	ng	# 0.02
29) Chrysene-d12	21.304	240	58869	0.400	ng	0.02
35) Perylene-d12	23.599	264	41358	0.400	ng	0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.345	112	11724	0.247	ng	0.02
5) Phenol-d6	6.912	99	8176	0.182	ng	0.00
8) Nitrobenzene-d5	8.859	82	13559	0.291	ng	0.00
11) 2-Methylnaphthalene-d10	12.107	152	34820	0.249	ng	0.02
14) 2,4,6-Tribromophenol	15.895	330	880m	0.038	ng	0.05
15) 2-Fluorobiphenyl	12.990	172	39899m	0.246	ng	0.02
27) Fluoranthene-d10	19.139	212	85390	0.331	ng	0.00
31) Terphenyl-d14	19.750	244	64577	0.511	ng	0.01

Target Compounds Qvalue

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

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