

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100923\
 Data File : BN028174.D
 Acq On : 10 Oct 2023 00:14
 Operator : MA/JU
 Sample : PB156183BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB156183BL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/10/2023
 Supervised By :mohammad ahmed 10/10/2023

Quant Time: Oct 10 00:58:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 10 00:42:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3880	0.400	ng	0.00
7) Naphthalene-d8	10.746	136	10652	0.400	ng	0.01
13) Acenaphthene-d10	14.576	164	5365	0.400	ng	0.01
19) Phenanthrene-d10	17.331	188	10259	0.400	ng	# 0.01
29) Chrysene-d12	21.515	240	8929	0.400	ng	# 0.00
35) Perylene-d12	23.978	264	9190	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.466	112	4128	0.357	ng	0.00
5) Phenol-d6	7.098	99	4072	0.279	ng	0.00
8) Nitrobenzene-d5	9.144	82	2945	0.325	ng	0.02
11) 2-Methylnaphthalene-d10	12.347	152	5901	0.374	ng	0.02
14) 2,4,6-Tribromophenol	16.090	330	605	0.246	ng	0.01
15) 2-Fluorobiphenyl	13.218	172	7473m	0.387	ng	0.03
27) Fluoranthene-d10	19.360	212	10456	0.406	ng	0.00
31) Terphenyl-d14	19.946	244	8542	0.410	ng	0.00

Target Compounds Qvalue

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

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