

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034558.D
 Acq On : 11 Oct 2024 16:59
 Operator : RC/JU
 Sample : PB164072BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB164072BL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/14/2024
 Supervised By :mohammad ahmed 10/15/2024

Quant Time: Oct 11 17:28:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	4651	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	12810	0.400	ng	0.01
13) Acenaphthene-d10	14.030	164	7198	0.400	ng	0.01
19) Phenanthrene-d10	16.796	188	11872	0.400	ng	# 0.02
29) Chrysene-d12	21.008	240	7454	0.400	ng	0.00
35) Perylene-d12	23.110	264	10379	0.400	ng	# 0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	3274	0.248	ng	0.00
5) Phenol-d6	6.629	99	2846m	0.185	ng	0.06
8) Nitrobenzene-d5	8.546	82	2145m	0.219	ng	0.03
11) 2-Methylnaphthalene-d10	11.761	152	4835m	0.256	ng	0.03
14) 2,4,6-Tribromophenol	15.567	330	459	0.141	ng	0.03
15) 2-Fluorobiphenyl	12.661	172	5819	0.199	ng	0.03
27) Fluoranthene-d10	18.834	212	8217	0.274	ng	0.00
31) Terphenyl-d14	19.452	244	5159	0.347	ng	0.00

Target Compounds Qvalue

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

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