

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122324\
 Data File : BN035831.D
 Acq On : 24 Dec 2024 22:07
 Operator : RC/JU
 Sample : PB165596BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB165596BL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/25/2024
 Supervised By :mohammad ahmed 12/26/2024

Quant Time: Dec 24 23:55:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN112724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 19 23:06:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.264	152	4175	0.400	ng	0.00
7) Naphthalene-d8	10.009	136	9064	0.400	ng	-0.01
13) Acenaphthene-d10	13.935	164	5565	0.400	ng	0.00
19) Phenanthrene-d10	16.711	188	10427	0.400	ng	# 0.01
29) Chrysene-d12	20.956	240	8385	0.400	ng	# 0.00
35) Perylene-d12	23.035	264	8128	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.924	112	3132	0.300	ng	-0.01
5) Phenol-d6	6.477	99	2766	0.220	ng	-0.01
8) Nitrobenzene-d5	8.418	82	2502m	0.452	ng	0.01
11) 2-Methylnaphthalene-d10	11.636	152	4854	0.342	ng	0.02
14) 2,4,6-Tribromophenol	15.475	330	385	0.097	ng	0.02
15) 2-Fluorobiphenyl	12.559	172	7363	0.350	ng	0.02
27) Fluoranthene-d10	18.761	212	11459	0.388	ng	0.00
31) Terphenyl-d14	19.388	244	9065	0.548	ng	0.00

Target Compounds Qvalue

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

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