

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121924\
 Data File : BN035713.D
 Acq On : 19 Dec 2024 20:35
 Operator : RC/JU
 Sample : PB165596BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB165596BL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/21/2024
 Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 19 23:11:28 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	7.272	152	3228	0.400	ng	0.00
7) Naphthalene-d8	10.020	136	8561	0.400	ng	0.00
13) Acenaphthene-d10	13.935	164	5768	0.400	ng	0.00
19) Phenanthrene-d10	16.711	188	10744	0.400	ng	# 0.01
29) Chrysene-d12	20.956	240	9839	0.400	ng	# 0.00
35) Perylene-d12	23.044	264	11233	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.939	112	2902	0.359	ng	0.00
5) Phenol-d6	6.492	99	2957	0.304	ng	0.00
8) Nitrobenzene-d5	8.418	82	2280m	0.436	ng	0.01
11) 2-Methylnaphthalene-d10	11.631	152	5143	0.384	ng	0.00
14) 2,4,6-Tribromophenol	15.464	330	464	0.113	ng	0.01
15) 2-Fluorobiphenyl	12.549	172	7059	0.324	ng	0.00
27) Fluoranthene-d10	18.766	212	11890	0.390	ng	0.00
31) Terphenyl-d14	19.393	244	10159	0.523	ng	0.00

Target Compounds Qvalue

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

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