Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122324\
Data File : BN035808.D
Acq On : 24 Dec 2024 06:34
Operator : RC/JU
Sample : PB165667BS
Misc :
ALS Vial : 92 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PB165667BS

Manual Integrations APPROVED
Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024

Quant Time: Dec 24 07:07:05 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	4067	0.400	ng	0.00
7) Naphthalene-d8	10.009	136	9754	0.400	ng	#-0.01
13) Acenaphthene-d10	13.914	164	5527	0.400	ng	-0.02
19) Phenanthrene-d10	16.698	188	10920	0.400	ng	# 0.00
29) Chrysene-d12	20.947	240	9467	0.400	ng	# 0.00
35) Perylene-d12	23.032	264	9435	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.932	112	3134	0.308	ng	0.00
5) Phenol-d6	6.477	99	3400	0.278	ng	-0.01
8) Nitrobenzene-d5	8.397	82	2771m	0.465	ng	-0.01
11) 2-Methylnaphthalene-d10	11.611	152	7007	0.459	ng	-0.01
14) 2,4,6-Tribromophenol	15.453	330	647	0.165	ng	0.00
15) 2-Fluorobiphenyl	12.528	172	8330	0.399	ng	-0.02
27) Fluoranthene-d10	18.752	212	11348	0.367	ng	0.00
31) Terphenyl-d14	19.384	244	8056	0.431	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.982	88	1441	0.371	ng	# 72
3) n-Nitrosodimethylamine	3.271	42	1276	0.394	ng	# 88
6) bis(2-Chloroethyl)ether	6.715	93	3429	0.333	ng	99
9) Naphthalene	10.052	128	9192	0.357	ng	100
10) Hexachlorobutadiene	10.340	225	2323	0.391	ng	# 99
12) 2-Methylnaphthalene	11.687	142	5969	0.324	ng	98
16) Acenaphthylene	13.636	152	8595	0.370	ng	99
17) Acenaphthene	13.978	154	5618	0.365	ng	99
18) Fluorene	14.994	166	7068	0.320	ng	99
20) 4,6-Dinitro-2-methylph...	15.122	198	260	0.242	ng	# 30
21) 4-Bromophenyl-phenylether	15.904	248	2229	0.349	ng	# 91
22) Hexachlorobenzene	16.003	284	3165	0.422	ng	98
23) Atrazine	16.202	200	1619	0.356	ng	96
24) Pentachlorophenol	16.363	266	1112	0.341	ng	# 84
25) Phenanthrene	16.736	178	10790	0.360	ng	99
26) Anthracene	16.835	178	9628	0.355	ng	99
28) Fluoranthene	18.785	202	12957	0.320	ng	100
30) Pyrene	19.152	202	13632	0.390	ng	99
32) Benzo(a)anthracene	20.929	228	10627	0.321	ng	99
33) Chrysene	20.983	228	14185	0.415	ng	100
34) Bis(2-ethylhexyl)phtha...	20.902	149	4640	0.355	ng	100
36) Indeno(1,2,3-cd)pyrene	25.053	276	13497	0.366	ng	98
37) Benzo(b)fluoranthene	22.421	252	22672	0.657	ng	96
38) Benzo(k)fluoranthene	22.462	252	14330	0.422	ng	97
39) Benzo(a)pyrene	22.942	252	11221	0.395	ng	# 92
40) Dibenzo(a,h)anthracene	25.076	278	9782	0.336	ng	93
41) Benzo(g,h,i)perylene	25.664	276	11572	0.380	ng	99

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

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122324\
Data File : BN035808.D
Acq On : 24 Dec 2024 06:34
Operator : RC/JU
Sample : PB165667BS
Misc :
ALS Vial : 92 Sample Multiplier: 1

Quant Time: Dec 24 07:07:05 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

Instrument :
BNA_N
ClientSampleId :
PB165667BS

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024