Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121924\
Data File : BN035764.D
Acq On : 21 Dec 2024 04:34
Operator : RC/JU
Sample : PB165724BS
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PB165724BS

Manual Integrations APPROVED
Reviewed By :Jagrut Upadhyay 12/23/2024
Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 21 05:09:33 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	7.271	152	2647	0.400	ng	0.00
7) Naphthalene-d8	10.009	136	6538	0.400	ng	-0.01
13) Acenaphthene-d10	13.924	164	3740	0.400	ng	-0.01
19) Phenanthrene-d10	16.698	188	6891	0.400	ng	# 0.00
29) Chrysene-d12	20.955	240	6144	0.400	ng	0.00
35) Perylene-d12	23.041	264	6576	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.939	112	2510	0.379	ng	0.00
5) Phenol-d6	6.484	99	2665	0.334	ng	0.00
8) Nitrobenzene-d5	8.407	82	2003m	0.502	ng	0.00
11) 2-Methylnaphthalene-d10	11.615	152	5689	0.556	ng	0.00
14) 2,4,6-Tribromophenol	15.453	330	453	0.171	ng	0.00
15) 2-Fluorobiphenyl	12.538	172	6506	0.460	ng	0.00
27) Fluoranthene-d10	18.761	212	8504	0.435	ng	0.00
31) Terphenyl-d14	19.393	244	6077	0.501	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.981	88	1209	0.478	ng	# 64
3) n-Nitrosodimethylamine	3.270	42	1089	0.517	ng	# 97
6) bis(2-Chloroethyl)ether	6.722	93	2779	0.415	ng	99
9) Naphthalene	10.062	128	7464	0.433	ng	99
10) Hexachlorobutadiene	10.351	225	1928	0.485	ng	# 99
12) 2-Methylnaphthalene	11.691	142	4846	0.393	ng	97
16) Acenaphthylene	13.636	152	6779	0.432	ng	99
17) Acenaphthene	13.988	154	4529	0.434	ng	98
18) Fluorene	14.993	166	5492	0.368	ng	100
20) 4,6-Dinitro-2-methylph...	15.143	198	74	0.109	ng	# 1
21) 4-Bromophenyl-phenylether	15.916	248	1621	0.402	ng	# 92
22) Hexachlorobenzene	16.003	284	2441	0.516	ng	96
23) Atrazine	16.214	200	1272	0.443	ng	94
24) Pentachlorophenol	16.375	266	1192	0.579	ng	# 84
25) Phenanthrene	16.748	178	7957	0.420	ng	99
26) Anthracene	16.835	178	7046	0.412	ng	100
28) Fluoranthene	18.794	202	9671	0.379	ng	99
30) Pyrene	19.161	202	10032	0.442	ng	99
32) Benzo(a)anthracene	20.937	228	8148	0.379	ng	99
33) Chrysene	20.991	228	10584	0.478	ng	99
34) Bis(2-ethylhexyl)phtha...	20.902	149	3563m	0.420	ng	
36) Indeno(1,2,3-cd)pyrene	25.067	276	11390	0.443	ng	96
37) Benzo(b)fluoranthene	22.430	252	13899	0.578	ng	# 94
38) Benzo(k)fluoranthene	22.471	252	10950	0.462	ng	94
39) Benzo(a)pyrene	22.956	252	8888	0.449	ng	# 88
40) Dibenzo(a,h)anthracene	25.090	278	8351	0.412	ng	95
41) Benzo(g,h,i)perylene	25.678	276	9596	0.453	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121924\
Data File : BN035764.D
Acq On : 21 Dec 2024 04:34
Operator : RC/JU
Sample : PB165724BS
Misc :
ALS Vial : 55 Sample Multiplier: 1

Quant Time: Dec 21 05:09:33 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 :
PB165724BS

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/23/2024
Supervised By :mohammad ahmed 12/24/2024