Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040325\
Data File : BN036827.D
Acq On : 03 Apr 2025 17:50
Operator : RC/JU
Sample : PB167430BSD
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Apr 03 18:14:02 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN031025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 10 16:06:28 2025
Response via : Initial Calibration

Instrument :
BNA_N
ClientSampleId :
PB167430BSD

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 04/04/2025
Supervised By :Jagrut Upadhyay 04/04/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.695	152	2152	0.400	ng	-0.03
7) Naphthalene-d8	10.477	136	5190	0.400	ng	-0.03
13) Acenaphthene-d10	14.334	164	2805	0.400	ng	-0.03
19) Phenanthrene-d10	17.086	188	5710	0.400	ng	-0.02
29) Chrysene-d12	21.277	240	4059	0.400	ng	#-0.02
35) Perylene-d12	23.516	264	1954	0.400	ng	#-0.04
System Monitoring Compounds						
4) 2-Fluorophenol	5.290	112	2032	0.405	ng	-0.02
5) Phenol-d6	6.872	99	2364	0.382	ng	-0.03
8) Nitrobenzene-d5	8.843	82	1762	0.312	ng	-0.03
11) 2-Methylnaphthalene-d10	12.070	152	3326m	0.431	ng	-0.04
14) 2,4,6-Tribromophenol	15.833	330	465	0.365	ng	-0.02
15) 2-Fluorobiphenyl	12.958	172	5694	0.349	ng	-0.03
27) Fluoranthene-d10	19.113	212	5064	0.346	ng	-0.03
31) Terphenyl-d14	19.717	244	3498	0.360	ng	-0.03
Target Compounds						Qvalue
2) 1,4-Dioxane	3.225	88	841	0.352	ng	# 38
3) n-Nitrosodimethylamine	3.528	42	1924	0.398	ng	# 95
6) bis(2-Chloroethyl)ether	7.118	93	2319	0.362	ng	97
9) Naphthalene	10.530	128	5604	0.367	ng	100
10) Hexachlorobutadiene	10.818	225	1249	0.348	ng	# 100
12) 2-Methylnaphthalene	12.146	142	3541	0.365	ng	99
16) Acenaphthylene	14.056	152	5028	0.380	ng	99
17) Acenaphthene	14.398	154	3346	0.386	ng	99
18) Fluorene	15.382	166	4448	0.379	ng	99
20) 4,6-Dinitro-2-methylph...	15.478	198	472	0.469	ng	# 80
21) 4-Bromophenyl-phenylether	16.280	248	1317	0.368	ng	93
22) Hexachlorobenzene	16.391	284	1540	0.357	ng	97
23) Atrazine	16.565	200	325	0.113	ng	# 84
24) Pentachlorophenol	16.739	266	1406	0.714	ng	99
25) Phenanthrene	17.124	178	6736	0.393	ng	99
26) Anthracene	17.210	178	6105	0.395	ng	100
28) Fluoranthene	19.146	202	7399	0.385	ng	99
30) Pyrene	19.508	202	6876	0.346	ng	100
32) Benzo(a)anthracene	21.259	228	5364	0.380	ng	100
33) Chrysene	21.313	228	6613	0.429	ng	98
34) Bis(2-ethylhexyl)phtha...	21.196	149	3514	0.350	ng	# 99
36) Indeno(1,2,3-cd)pyrene	25.782	276	4405	0.625	ng	99
37) Benzo(b)fluoranthene	22.841	252	5131	0.722	ng	97
38) Benzo(k)fluoranthene	22.888	252	5174	0.694	ng	96
39) Benzo(a)pyrene	23.420	252	3260	0.544	ng	95
40) Dibenzo(a,h)anthracene	25.803	278	4093	0.745	ng	96
41) Benzo(g,h,i)perylene	26.466	276	3441	0.548	ng	99

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

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