Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013025\
Data File : BF141315.D
Acq On : 30 Jan 2025 16:03
Operator : RC/JU
Sample : Q1209-05MS
Misc :
ALS Vial : 16 Sample Multiplier: 1

Quant Time: Jan 30 16:35:20 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 29 17:41:25 2025
Response via : Initial Calibration

Instrument :
BNA_F
ClientSampleId :
WC-5MS

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay

01/30/2025
Supervised By :mohammad ahmed

01/31/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	167075	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	669898	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	361408	20.000	ng	0.00
64) Phenanthrene-d10	11.328	188	589870	20.000	ng	0.00
76) Chrysene-d12	13.974	240	319317	20.000	ng	0.00
86) Perylene-d12	15.445	264	315915	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	958938	88.114	ng	0.00
7) Phenol-d6	6.428	99	1223169	88.351	ng	-0.01
23) Nitrobenzene-d5	7.363	82	782312	61.563	ng	0.00
42) 2,4,6-Tribromophenol	10.628	330	309128	93.300	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1347404	57.384	ng	0.00
79) Terphenyl-d14	12.910	244	1349280	65.165	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.610	88	183718	38.384	ng	100
3) Pyridine	3.328	79	429489	37.243	ng	99
4) n-Nitrosodimethylamine	3.287	42	258831	39.289	ng	100
6) Aniline	6.463	93	392868	33.523	ng	98
8) 2-Chlorophenol	6.581	128	457640	39.233	ng	98
9) Benzaldehyde	6.351	77	407982	50.876	ng	98
10) Phenol	6.445	94	572177	38.988	ng	95
11) bis(2-Chloroethyl)ether	6.534	93	435282	38.683	ng	100
12) 1,3-Dichlorobenzene	6.740	146	466993	36.310	ng	99
13) 1,4-Dichlorobenzene	6.816	146	475503	36.802	ng	100
14) 1,2-Dichlorobenzene	6.969	146	455246	37.628	ng	98
15) Benzyl Alcohol	6.940	79	388911	41.106	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.075	45	789105	40.328	ng	99
17) 2-Methylphenol	7.051	107	373745	39.424	ng	98
18) Hexachloroethane	7.310	117	176604	37.051	ng	98
19) n-Nitroso-di-n-propyla...	7.210	70	309827	37.662	ng	100
20) 3+4-Methylphenols	7.204	107	434424	37.321	ng	92
22) Acetophenone	7.210	105	654231	41.086	ng	98
24) Nitrobenzene	7.381	77	481987	36.600	ng	99
25) Isophorone	7.622	82	869191	39.569	ng	99
26) 2-Nitrophenol	7.698	139	244478	39.364	ng	98
27) 2,4-Dimethylphenol	7.734	122	388605	49.159	ng	99
28) bis(2-Chloroethoxy)met...	7.834	93	542293	38.703	ng	99
29) 2,4-Dichlorophenol	7.940	162	373371	38.581	ng	99
30) 1,2,4-Trichlorobenzene	8.022	180	394571	35.701	ng	99
31) Naphthalene	8.104	128	1303056	36.821	ng	100
32) Benzoic acid	7.857	122	305768	42.076	ng	99
33) 4-Chloroaniline	8.151	127	180833	15.069	ng	99
34) Hexachlorobutadiene	8.222	225	234884	36.240	ng	99
35) Caprolactam	8.522	113	135885m	42.993	ng	
36) 4-Chloro-3-methylphenol	8.634	107	409952	39.492	ng	97
37) 2-Methylnaphthalene	8.792	142	818250	36.379	ng	100
38) 1-Methylnaphthalene	8.892	142	799890	36.203	ng	100
40) 1,2,4,5-Tetrachloroben...	8.957	216	416191	40.479	ng	99
41) Hexachlorocyclopentadiene	8.945	237	449360	147.988	ng	100
43) 2,4,6-Trichlorophenol	9.069	196	271365	40.323	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	262590	35.851	ng	99
46) 1,1'-Biphenyl	9.257	154	1148416	40.745	ng	99
47) 2-Chloronaphthalene	9.281	162	796174	36.241	ng	99
48) 2-Nitroaniline	9.381	65	268545	38.862	ng	99
49) Acenaphthylene	9.698	152	1236080	40.194	ng	100
50) Dimethylphthalate	9.563	163	931376	38.156	ng	99
51) 2,6-Dinitrotoluene	9.622	165	206068	36.359	ng	100
52) Acenaphthene	9.869	154	910445	41.441	ng	100
53) 3-Nitroaniline	9.786	138	99441	17.902	ng	97
54) 2,4-Dinitrophenol	9.898	184	245895	93.333	ng	97
55) Dibenzofuran	10.045	168	1093675	37.190	ng	99
56) 4-Nitrophenol	9.957	139	360991	88.860	ng	98
57) 2,4-Dinitrotoluene	10.028	165	280902	38.281	ng	96
58) Fluorene	10.386	166	840449	36.844	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.163	232	231917	40.247	ng	99
60) Diethylphthalate	10.263	149	893576	37.755	ng	100
61) 4-Chlorophenyl-phenyle...	10.381	204	405320	36.376	ng	99
62) 4-Nitroaniline	10.410	138	211996	38.819	ng	99
63) Azobenzene	10.539	77	1037084	43.776	ng	98
65) 4,6-Dinitro-2-methylph...	10.433	198	156125	44.747	ng	96
66) n-Nitrosodiphenylamine	10.498	169	757543	39.707	ng	99
67) 4-Bromophenyl-phenylether	10.869	248	258485	39.142	ng	98
68) Hexachlorobenzene	10.933	284	268678	38.355	ng	100
69) Atrazine	11.028	200	276513	43.348	ng	99
70) Pentachlorophenol	11.128	266	338623	82.551	ng	99
71) Phenanthrene	11.351	178	1246279	39.305	ng	100
72) Anthracene	11.404	178	1270809	40.904	ng	100
73) Carbazole	11.557	167	1104522	40.093	ng	99
74) Di-n-butylphthalate	11.892	149	1342332	40.206	ng	100
75) Fluoranthene	12.539	202	1199766	38.321	ng	100
77) Benzidine	12.663	184	534831	52.246	ng	100
78) Pyrene	12.769	202	1190370	38.836	ng	100
80) Butylbenzylphthalate	13.392	149	473214	44.191	ng	100
81) Benzo(a)anthracene	13.963	228	812023	36.852	ng	100
82) 3,3'-Dichlorobenzidine	13.927	252	205049	29.259	ng	99
83) Chrysene	13.998	228	784739	38.860	ng	100
84) Bis(2-ethylhexyl)phtha...	13.957	149	575904	42.396	ng	99
85) Di-n-octyl phthalate	14.592	149	845371	40.068	ng	100
87) Indeno(1,2,3-cd)pyrene	16.904	276	916478	44.947	ng	99
88) Benzo(b)fluoranthene	15.027	252	756397	38.064	ng	99
89) Benzo(k)fluoranthene	15.057	252	677523	38.487	ng	100
90) Benzo(a)pyrene	15.386	252	704442	41.949	ng	100
91) Dibenzo(a,h)anthracene	16.921	278	733439	44.710	ng	99
92) Benzo(g,h,i)perylene	17.339	276	711929	41.738	ng	99

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

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