Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013125\
Data File : BF141334.D
Acq On : 31 Jan 2025 16:26
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSTDCCC040EC

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 02/03/2025
Supervised By :mohammad ahmed 02/04/2025

Quant Time: Jan 31 18:34:55 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	164970	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	647909	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	341541	20.000	ng	0.00
64) Phenanthrene-d10	11.321	188	566479	20.000	ng	0.00
76) Chrysene-d12	13.968	240	341440	20.000	ng	0.00
86) Perylene-d12	15.439	264	331116	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	791139	73.623	ng	0.00
7) Phenol-d6	6.428	99	993362	72.667	ng	-0.01
23) Nitrobenzene-d5	7.363	82	896200	72.918	ng	0.00
42) 2,4,6-Tribromophenol	10.627	330	235886	75.335	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1562690	70.424	ng	0.00
79) Terphenyl-d14	12.915	244	1567460	70.798	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.481	88	179518	37.986	ng	99
3) Pyridine	3.210	79	436407	38.326	ng	99
4) n-Nitrosodimethylamine	3.163	42	248931	38.269	ng	99
6) Aniline	6.463	93	439932	38.018	ng	99
8) 2-Chlorophenol	6.586	128	426501	37.030	ng	96
9) Benzaldehyde	6.345	77	310932	39.268	ng	99
10) Phenol	6.445	94	527421	36.397	ng	94
11) bis(2-Chloroethyl)ether	6.539	93	407197	36.649	ng	98
12) 1,3-Dichlorobenzene	6.739	146	469100	36.939	ng	99
13) 1,4-Dichlorobenzene	6.816	146	473169	37.089	ng	100
14) 1,2-Dichlorobenzene	6.969	146	432488	36.203	ng	99
15) Benzyl Alcohol	6.939	79	350130	37.479	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.075	45	715839	37.050	ng	99
17) 2-Methylphenol	7.051	107	344483	36.801	ng	99
18) Hexachloroethane	7.310	117	173871	36.943	ng	98
19) n-Nitroso-di-n-propyla...	7.216	70	295405	36.368	ng	97
20) 3+4-Methylphenols	7.204	107	411022	35.761	ng	99
22) Acetophenone	7.210	105	548834	35.637	ng	99
24) Nitrobenzene	7.381	77	467777	36.727	ng	99
25) Isophorone	7.622	82	785054	36.952	ng	99
26) 2-Nitrophenol	7.698	139	224162	37.318	ng	100
27) 2,4-Dimethylphenol	7.733	122	286982m	37.535	ng	
28) bis(2-Chloroethoxy)met...	7.833	93	494480	36.489	ng	100
29) 2,4-Dichlorophenol	7.939	162	346397	37.008	ng	100
30) 1,2,4-Trichlorobenzene	8.022	180	385139	36.030	ng	99
31) Naphthalene	8.104	128	1231790	35.989	ng	99
32) Benzoic acid	7.851	122	298299m	42.441	ng	
33) 4-Chloroaniline	8.151	127	411817	35.481	ng	99
34) Hexachlorobutadiene	8.222	225	227599	36.308	ng	99
35) Caprolactam	8.528	113	116667	38.166	ng	94
36) 4-Chloro-3-methylphenol	8.633	107	376529	37.503	ng	98
37) 2-Methylnaphthalene	8.792	142	782487	35.969	ng	98
38) 1-Methylnaphthalene	8.892	142	767501	35.916	ng	100
40) 1,2,4,5-Tetrachloroben...	8.957	216	357735	36.817	ng	100
41) Hexachlorocyclopentadiene	8.945	237	112765	39.297	ng	99
43) 2,4,6-Trichlorophenol	9.069	196	241760	38.014	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	256970	37.124	ng	100
46) 1,1'-Biphenyl	9.257	154	961081	36.082	ng	99
47) 2-Chloronaphthalene	9.286	162	748367	36.046	ng	99
48) 2-Nitroaniline	9.380	65	246493	37.746	ng	99
49) Acenaphthylene	9.698	152	1061017	36.509	ng	100
50) Dimethylphthalate	9.563	163	853360	36.994	ng	100
51) 2,6-Dinitrotoluene	9.622	165	202525	37.813	ng	98
52) Acenaphthene	9.875	154	761516	36.678	ng	100
53) 3-Nitroaniline	9.792	138	195318	37.207	ng	100
54) 2,4-Dinitrophenol	9.898	184	104461	41.956	ng	97
55) Dibenzofuran	10.045	168	999283	35.957	ng	100
56) 4-Nitrophenol	9.951	139	152830	39.808	ng	98
57) 2,4-Dinitrotoluene	10.027	165	267112	38.519	ng	99
58) Fluorene	10.386	166	773726	35.892	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.163	232	209383	38.450	ng	99
60) Diethylphthalate	10.263	149	832152	37.205	ng	100
61) 4-Chlorophenyl-phenyle...	10.380	204	380628	36.147	ng	100
62) 4-Nitroaniline	10.404	138	198535	38.469	ng	99
63) Azobenzene	10.539	77	831958	37.160	ng	99
65) 4,6-Dinitro-2-methylph...	10.433	198	148537	44.330	ng	99
66) n-Nitrosodiphenylamine	10.498	169	682549	37.253	ng	99
67) 4-Bromophenyl-phenylether	10.869	248	244392	38.536	ng	98
68) Hexachlorobenzene	10.933	284	255323	37.954	ng	99
69) Atrazine	11.027	200	213403	34.836	ng	98
70) Pentachlorophenol	11.127	266	167246	42.456	ng	99
71) Phenanthrene	11.351	178	1133127	37.212	ng	99
72) Anthracene	11.404	178	1126615	37.760	ng	99
73) Carbazole	11.557	167	1014712	38.354	ng	99
74) Di-n-butylphthalate	11.892	149	1270081	39.613	ng	100
75) Fluoranthene	12.539	202	1159837	38.576	ng	100
77) Benzidine	12.663	184	393366	35.937	ng	99
78) Pyrene	12.768	202	1173910	35.817	ng	100
80) Butylbenzylphthalate	13.392	149	451201	39.406	ng	99
81) Benzo(a)anthracene	13.957	228	824181	34.980	ng	100
82) 3,3'-Dichlorobenzidine	13.927	252	272114	36.313	ng	99
83) Chrysene	13.998	228	770666	35.690	ng	100
84) Bis(2-ethylhexyl)phtha...	13.957	149	561147	38.633	ng	100
85) Di-n-octyl phthalate	14.586	149	880674	39.037	ng	100
87) Indeno(1,2,3-cd)pyrene	16.892	276	802001	37.527	ng	100
88) Benzo(b)fluoranthene	15.015	252	808478	38.817	ng	99
89) Benzo(k)fluoranthene	15.045	252	638999	34.632	ng	100
90) Benzo(a)pyrene	15.380	252	655984	37.270	ng	100
91) Dibenzo(a,h)anthracene	16.909	278	645395	37.536	ng	99
92) Benzo(g,h,i)perylene	17.327	276	685709	38.355	ng	99

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

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