

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121223\
 Data File : BF136508.D
 Acq On : 12 Dec 2023 10:52
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 12 11:40:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 06 09:40:27 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/14/2023
 Supervised By :mohammad ahmed 12/14/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	118805	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	487987	20.000	ng	0.00
39) Acenaphthene-d10	9.833	164	262692	20.000	ng	0.00
64) Phenanthrene-d10	11.328	188	495404	20.000	ng	0.00
76) Chrysene-d12	13.980	240	254997	20.000	ng	0.00
86) Perylene-d12	15.457	264	246710	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	595419	74.015	ng	0.00
7) Phenol-d6	6.445	99	748486	74.573	ng	0.00
23) Nitrobenzene-d5	7.363	82	695628	77.063	ng	0.00
42) 2,4,6-Tribromophenol	10.628	330	248673	76.774	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1439156	74.754	ng	0.00
79) Terphenyl-d14	12.922	244	1591590	82.267	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.557	88	116901	36.740	ng	98
3) Pyridine	3.316	79	316771	41.109	ng	95
4) n-Nitrosodimethylamine	3.287	42	126278	37.720	ng	95
6) Aniline	6.469	93	467537	46.093	ng	99
8) 2-Chlorophenol	6.587	128	335817	38.238	ng	97
9) Benzaldehyde	6.357	77	208994	41.575	ng	96
10) Phenol	6.457	94	402565	37.693	ng	90
11) bis(2-Chloroethyl)ether	6.540	93	301408	37.861	ng	99
12) 1,3-Dichlorobenzene	6.740	146	346059	35.038	ng	99
13) 1,4-Dichlorobenzene	6.816	146	351680	35.101	ng	99
14) 1,2-Dichlorobenzene	6.969	146	327080	34.717	ng	99
15) Benzyl Alcohol	6.945	79	271163	40.346	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.075	45	370822	36.217	ng	96
17) 2-Methylphenol	7.057	107	284340	40.075	ng	97
18) Hexachloroethane	7.304	117	121779	34.814	ng	96
19) n-Nitroso-di-n-propyla...	7.216	70	222286	38.393	ng	99
20) 3+4-Methylphenols	7.210	107	354443	40.246	ng	95
22) Acetophenone	7.210	105	458994	38.090	ng	98
24) Nitrobenzene	7.381	77	334226	36.806	ng	99
25) Isophorone	7.622	82	630322	38.867	ng	98
26) 2-Nitrophenol	7.698	139	181211	40.592	ng	91
27) 2,4-Dimethylphenol	7.739	122	276853	50.294	ng	100
28) bis(2-Chloroethoxy)met...	7.834	93	389920	39.128	ng	99
29) 2,4-Dichlorophenol	7.939	162	283869	39.069	ng	100
30) 1,2,4-Trichlorobenzene	8.016	180	309893	36.577	ng	98
31) Naphthalene	8.098	128	999471	37.141	ng	99
32) Benzoic acid	7.869	122	212136m	38.848	ng	
33) 4-Chloroaniline	8.151	127	420567	43.837	ng	99
34) Hexachlorobutadiene	8.216	225	182476	36.342	ng	98
35) Caprolactam	8.534	113	91946	41.093	ng	95
36) 4-Chloro-3-methylphenol	8.639	107	306083	39.675	ng	98
37) 2-Methylnaphthalene	8.792	142	686358	40.091	ng	96
38) 1-Methylnaphthalene	8.892	142	639020	38.038	ng	98
40) 1,2,4,5-Tetrachloroben...	8.957	216	316832	38.767	ng	100
41) Hexachlorocyclopentadiene	8.939	237	135396	44.718	ng	98
43) 2,4,6-Trichlorophenol	9.069	196	204618	38.451	ng	97

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44) 2,4,5-Trichlorophenol	9.116	196	238119	40.092	ng	93
46) 1,1'-Biphenyl	9.257	154	841664	37.654	ng	99
47) 2-Chloronaphthalene	9.281	162	628351	36.118	ng	97
48) 2-Nitroaniline	9.381	65	186512	39.972	ng	97
49) Acenaphthylene	9.698	152	976683	39.409	ng	99
50) Dimethylphthalate	9.563	163	773068	39.972	ng	98
51) 2,6-Dinitrotoluene	9.628	165	169932	39.056	ng	94
52) Acenaphthene	9.869	154	602688	37.950	ng	98
53) 3-Nitroaniline	9.798	138	181959	40.953	ng	97
54) 2,4-Dinitrophenol	9.904	184	85685	38.950	ng	96
55) Dibenzofuran	10.045	168	902120	38.884	ng	97
56) 4-Nitrophenol	9.963	139	125893	37.797	ng	92
57) 2,4-Dinitrotoluene	10.033	165	221729	38.522	ng	97
58) Fluorene	10.386	166	709893	38.118	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.163	232	182706	38.404	ng	99
60) Diethylphthalate	10.263	149	744622	39.154	ng	99
61) 4-Chlorophenyl-phenyle...	10.381	204	354776	39.077	ng	93
62) 4-Nitroaniline	10.410	138	179055	40.585	ng	95
63) Azobenzene	10.539	77	661574	38.219	ng	98
65) 4,6-Dinitro-2-methylph...	10.439	198	122937	40.069	ng	99
66) n-Nitrosodiphenylamine	10.498	169	637436	40.099	ng	98
67) 4-Bromophenyl-phenylether	10.869	248	226170	39.267	ng	93
68) Hexachlorobenzene	10.933	284	242541	36.433	ng	97
69) Atrazine	11.033	200	169700	37.403	ng	98
70) Pentachlorophenol	11.133	266	136386	36.300	ng	98
71) Phenanthrene	11.351	178	1061980	39.001	ng	99
72) Anthracene	11.404	178	1068116	38.945	ng	99
73) Carbazole	11.563	167	899956	37.075	ng	99
74) Di-n-butylphthalate	11.892	149	1139250	38.284	ng	99
75) Fluoranthene	12.545	202	1035589	36.292	ng	98
77) Benzidine	12.669	184	159572	31.777	ng	98
78) Pyrene	12.774	202	1022352	40.938	ng	98
80) Butylbenzylphthalate	13.398	149	367283	41.314	ng	91
81) Benzo(a)anthracene	13.969	228	692696	38.998	ng	99
82) 3,3'-Dichlorobenzidine	13.933	252	223842	44.947	ng	100
83) Chrysene	14.004	228	681597	39.003	ng	98
84) Bis(2-ethylhexyl)phtha...	13.963	149	442282	40.582	ng	99
85) Di-n-octyl phthalate	14.580	149	646156	40.139	ng	99
87) Indeno(1,2,3-cd)pyrene	16.962	276	706206	41.203	ng	99
88) Benzo(b)fluoranthene	15.027	252	580441	34.410	ng	99
89) Benzo(k)fluoranthene	15.057	252	562159	35.363	ng	99
90) Benzo(a)pyrene	15.398	252	552985	39.696	ng	99
91) Dibenzo(a,h)anthracene	16.986	278	580575	41.348	ng	99
92) Benzo(g,h,i)perylene	17.427	276	582168	40.395	ng	99

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

