

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
 Data File : BF136669.D
 Acq On : 21 Dec 2023 14:18
 Operator : CG\JU
 Sample : 05602-05MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SPRING-VALLEY-TP-3MSD

Quant Time: Dec 21 15:00:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/22/2023
 Supervised By :mohammad ahmed 12/22/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	72656	20.000	ng	-0.01
21) Naphthalene-d8	8.063	136	255830	20.000	ng	# 0.00
39) Acenaphthene-d10	9.816	164	103081	20.000	ng	-0.01
64) Phenanthrene-d10	11.304	188	177016	20.000	ng	-0.01
76) Chrysene-d12	13.963	240	145976	20.000	ng	0.00
86) Perylene-d12	15.439	264	122263	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	450360	96.715	ng	0.00
7) Phenol-d6	6.434	99	552575	91.580	ng	0.00
23) Nitrobenzene-d5	7.345	82	341224	68.252	ng	-0.01
42) 2,4,6-Tribromophenol	10.610	330	106874	88.034	ng	-0.01
45) 2-Fluorobiphenyl	9.139	172	580862	75.790	ng	-0.01
79) Terphenyl-d14	12.904	244	548074	52.469	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.575	88	74360	38.326	ng	99
3) Pyridine	3.334	79	179992	38.022	ng	98
4) n-Nitrosodimethylamine	3.305	42	116094	45.923	ng	98
6) Aniline	6.457	93	143699	23.820	ng	# 84
8) 2-Chlorophenol	6.575	128	227714	44.686	ng	95
9) Benzaldehyde	6.334	77	87165	25.398	ng	99
10) Phenol	6.445	94	277554	43.091	ng	92
11) bis(2-Chloroethyl)ether	6.522	93	279681	57.096	ng	97
12) 1,3-Dichlorobenzene	6.722	146	233748	42.048	ng	99
13) 1,4-Dichlorobenzene	6.798	146	239964	42.247	ng	99
14) 1,2-Dichlorobenzene	6.951	146	221639	41.934	ng	99
15) Benzyl Alcohol	6.934	79	181184	44.510	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.063	45	351909	43.942	ng	99
17) 2-Methylphenol	7.045	107	185357	43.907	ng	99
18) Hexachloroethane	7.287	117	84407	40.918	ng	100
19) n-Nitroso-di-n-propyla...	7.198	70	155368	42.477	ng	98
20) 3+4-Methylphenols	7.198	107	218957	41.516	ng	# 82
22) Acetophenone	7.192	105	311204	48.532	ng	# 94
24) Nitrobenzene	7.369	77	230743	46.073	ng	93
25) Isophorone	7.604	82	411783	45.265	ng	99
26) 2-Nitrophenol	7.681	139	114530	47.798	ng	96
27) 2,4-Dimethylphenol	7.722	122	164319	56.328	ng	98
28) bis(2-Chloroethoxy)met...	7.816	93	257434	46.555	ng	99
29) 2,4-Dichlorophenol	7.928	162	166248	44.267	ng	96
30) 1,2,4-Trichlorobenzene	8.004	180	183127	43.764	ng	98
31) Naphthalene	8.081	128	610608	43.445	ng	100
32) Benzoic acid	7.857	122	127593	45.205	ng	97
33) 4-Chloroaniline	8.145	127	144697	27.884	ng	95
34) Hexachlorobutadiene	8.198	225	111074	43.690	ng	97
35) Caprolactam	8.504	113	54740m	44.183	ng	
36) 4-Chloro-3-methylphenol	8.622	107	167989	39.733	ng	98
37) 2-Methylnaphthalene	8.775	142	361771	39.996	ng	99
38) 1-Methylnaphthalene	8.875	142	347267	39.133	ng	99
40) 1,2,4,5-Tetrachloroben...	8.939	216	171876	53.840	ng	98
41) Hexachlorocyclopentadiene	8.922	237	43362	44.891	ng	98
43) 2,4,6-Trichlorophenol	9.057	196	105949	51.750	ng	99

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44) 2,4,5-Trichlorophenol	9.098	196	109059	47.802	ng	95
46) 1,1'-Biphenyl	9.239	154	455578	52.668	ng	99
47) 2-Chloronaphthalene	9.263	162	323651	49.218	ng	99
48) 2-Nitroaniline	9.363	65	103081	50.058	ng	99
49) Acenaphthylene	9.681	152	500076	49.753	ng	100
50) Dimethylphthalate	9.545	163	376206	49.966	ng	99
51) 2,6-Dinitrotoluene	9.610	165	78013	45.654	ng	94
52) Acenaphthene	9.851	154	283002	45.422	ng	99
53) 3-Nitroaniline	9.775	138	67355	37.210	ng	97
54) 2,4-Dinitrophenol	9.886	184	72493	76.708	ng	# 84
55) Dibenzofuran	10.022	168	429155	45.439	ng	100
56) 4-Nitrophenol	9.951	139	106728	87.342	ng	97
57) 2,4-Dinitrotoluene	10.010	165	93442	42.123	ng	99
58) Fluorene	10.369	166	322708	43.062	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.145	232	79901	45.496	ng	99
60) Diethylphthalate	10.245	149	362689	46.427	ng	99
61) 4-Chlorophenyl-phenyle...	10.363	204	158170	43.423	ng	94
62) 4-Nitroaniline	10.392	138	65627m	37.608	ng	
63) Azobenzene	10.522	77	311458	42.686	ng	99
65) 4,6-Dinitro-2-methylph...	10.422	198	43736	43.112	ng	94
66) n-Nitrosodiphenylamine	10.481	169	285678	49.681	ng	99
67) 4-Bromophenyl-phenylether	10.851	248	90599	46.085	ng	98
68) Hexachlorobenzene	10.916	284	97885	44.680	ng	99
69) Atrazine	11.016	200	89535	60.877	ng	97
70) Pentachlorophenol	11.116	266	101105	99.010	ng	99
71) Phenanthrene	11.333	178	433203	47.695	ng	99
72) Anthracene	11.386	178	449998	48.921	ng	99
73) Carbazole	11.545	167	410132	47.291	ng	100
74) Di-n-butylphthalate	11.875	149	531267	50.202	ng	99
75) Fluoranthene	12.527	202	491722	50.612	ng	99
77) Benzidine	12.651	184	169030	39.271	ng	99
78) Pyrene	12.757	202	513557	35.838	ng	99
80) Butylbenzylphthalate	13.380	149	231602	44.417	ng	96
81) Benzo(a)anthracene	13.951	228	459405	45.322	ng	99
82) 3,3'-Dichlorobenzidine	13.916	252	155549	54.035	ng	96
83) Chrysene	13.992	228	452147	45.614	ng	99
84) Bis(2-ethylhexyl)phtha...	13.945	149	335929	51.205	ng	99
85) Di-n-octyl phthalate	14.563	149	614793	60.582	ng	99
87) Indeno(1,2,3-cd)pyrene	16.939	276	339234	38.429	ng	99
88) Benzo(b)fluoranthene	15.010	252	396944	48.609	ng	99
89) Benzo(k)fluoranthene	15.039	252	357942	46.389	ng	99
90) Benzo(a)pyrene	15.374	252	312648	45.357	ng	99
91) Dibenzo(a,h)anthracene	16.957	278	288595	39.849	ng	99
92) Benzo(g,h,i)perylene	17.392	276	262997	35.456	ng	98

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

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