

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062623\
 Data File : BF133969.D
 Acq On : 27 Jun 2023 01:59
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
ClientSampleId :
 SSTDCCC040

Manual Integrations
APPROVED
 Reviewed By :Yogesh Patel

Quant Time: Jun 27 06:02:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 23:12:04 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	93863	20.000	ng	0.00
21) Naphthalene-d8	8.134	136	300425	20.000	ng	0.00
39) Acenaphthene-d10	9.892	164	126292	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	257050	20.000	ng	0.00
76) Chrysene-d12	14.045	240	176767	20.000	ng	0.00
86) Perylene-d12	15.551	264	151132	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	472851	84.269	ng	0.00
7) Phenol-d6	6.493	99	531801	77.065	ng	0.00
23) Nitrobenzene-d5	7.416	82	475597	85.337	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	132476	83.663	ng	0.00
45) 2-Fluorobiphenyl	9.210	172	782425	90.074	ng	0.00
79) Terphenyl-d14	12.980	244	894120	81.407	ng	0.00

06/27/2023
 Supervised By :mohammad Ahmed
 06/29/2023

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.622	88	107476	46.452	ng	99
3) Pyridine	3.410	79	253214	42.016	ng	98
4) n-Nitrosodimethylamine	3.375	42	152724	44.371	ng	97
6) Aniline	6.522	93	328180	39.082	ng	100
8) 2-Chlorophenol	6.640	128	226276	39.725	ng	98
9) Benzaldehyde	6.410	77	153194	40.547	ng	97
10) Phenol	6.504	94	278537	38.389	ng	98
11) bis(2-Chloroethyl)ether	6.593	93	211135	39.526	ng	99
12) 1,3-Dichlorobenzene	6.793	146	246259	40.552	ng	99
13) 1,4-Dichlorobenzene	6.869	146	247852	40.408	ng	99
14) 1,2-Dichlorobenzene	7.022	146	224089	39.010	ng	98
15) Benzyl Alcohol	6.998	79	206525	38.822	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.128	45	364884	38.611	ng	98
17) 2-Methylphenol	7.104	107	188106	36.878	ng	98
18) Hexachloroethane	7.363	117	79038	34.752	ng	99
19) n-Nitroso-di-n-propyla...	7.263	70	158638	36.250	ng	95
20) 3+4-Methylphenols	7.257	107	226913	36.670	ng	# 86
22) Acetophenone	7.263	105	291321	42.638	ng	98
24) Nitrobenzene	7.434	77	239192	42.471	ng	94
25) Isophorone	7.669	82	408113	40.292	ng	97
26) 2-Nitrophenol	7.751	139	85554	31.667	ng	99
27) 2,4-Dimethylphenol	7.787	122	174214	40.737	ng	97
28) bis(2-Chloroethoxy)met...	7.881	93	242278	41.310	ng	99
29) 2,4-Dichlorophenol	7.992	162	160118	37.757	ng	97
30) 1,2,4-Trichlorobenzene	8.075	180	179401	40.999	ng	99
31) Naphthalene	8.157	128	582174	40.118	ng	99
32) Benzoic acid	7.863	122	68191m	21.279	ng	
33) 4-Chloroaniline	8.204	127	232206	37.408	ng	98
34) Hexachlorobutadiene	8.269	225	118704	43.005	ng	98
35) Caprolactam	8.569	113	45109	32.306	ng	92
36) 4-Chloro-3-methylphenol	8.687	107	168582	35.387	ng	98
37) 2-Methylnaphthalene	8.845	142	377217	36.759	ng	98
38) 1-Methylnaphthalene	8.945	142	345787	36.246	ng	100
40) 1,2,4,5-Tetrachloroben...	9.010	216	169708	45.357	ng	97
41) Hexachlorocyclopentadiene	8.992	237	19134	10.779	ng	93
43) 2,4,6-Trichlorophenol	9.128	196	108234	42.494	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.169	196	126097	42.088	ng	93	06/27/2023
46) 1,1'-Biphenyl	9.310	154	438891	43.323	ng	98	Supervised By :mohammad ahmed
47) 2-Chloronaphthalene	9.339	162	314571	42.439	ng	99	
48) 2-Nitroaniline	9.434	65	116051	42.960	ng	99	
49) Acenaphthylene	9.751	152	522261	41.246	ng	99	
50) Dimethylphthalate	9.610	163	389787	42.518	ng	99	
51) 2,6-Dinitrotoluene	9.675	165	76491	38.739	ng	87	06/29/2023
52) Acenaphthene	9.928	154	304286	41.415	ng	99	
53) 3-Nitroaniline	9.845	138	100303	42.343	ng	100	
54) 2,4-Dinitrophenol	9.957	184	2891	6.277	ng	# 80	
55) Dibenzofuran	10.098	168	473692	41.253	ng	100	
56) 4-Nitrophenol	10.010	139	60760	36.670	ng	96	
57) 2,4-Dinitrotoluene	10.081	165	96247	36.910	ng	93	
58) Fluorene	10.439	166	364246	40.774	ng	98	
59) 2,3,4,6-Tetrachlorophenol	10.216	232	96080	40.655	ng	96	
60) Diethylphthalate	10.310	149	400064	41.886	ng	98	
61) 4-Chlorophenyl-phenylether	10.434	204	176879	41.086	ng	97	
62) 4-Nitroaniline	10.463	138	100425	42.067	ng	96	
63) Azobenzene	10.598	77	367403	40.666	ng	99	
65) 4,6-Dinitro-2-methylphthalate	10.492	198	7445	5.312	ng	88	
66) n-Nitrosodiphenylamine	10.551	169	323134	39.345	ng	96	
67) 4-Bromophenyl-phenylether	10.928	248	108440	39.690	ng	99	
68) Hexachlorobenzene	10.992	284	115386	39.776	ng	99	
69) Atrazine	11.081	200	84365	37.137	ng	98	
70) Pentachlorophenol	11.186	266	57072	32.904	ng	97	
71) Phenanthrene	11.410	178	529216	40.897	ng	99	
72) Anthracene	11.463	178	551434	40.970	ng	99	
73) Carbazole	11.622	167	519924	42.203	ng	100	
74) Di-n-butylphthalate	11.951	149	635976	43.411	ng	99	
75) Fluoranthene	12.604	202	628307	44.912	ng	99	
77) Benzidine	12.733	184	208156	49.490	ng	98	
78) Pyrene	12.839	202	634891	38.594	ng	99	
80) Butylbenzylphthalate	13.463	149	291822	47.299	ng	95	
81) Benzo(a)anthracene	14.033	228	490838	40.039	ng	100	
82) 3,3'-Dichlorobenzidine	13.998	252	168539	43.394	ng	# 98	
83) Chrysene	14.074	228	467565	39.378	ng	99	
84) Bis(2-ethylhexyl)phthalate	14.022	149	392013	55.296	ng	98	
85) Di-n-octyl phthalate	14.645	149	600763	57.672	ng	100	
87) Indeno(1,2,3-cd)pyrene	17.098	276	356741	34.017	ng	98	
88) Benzo(b)fluoranthene	15.104	252	383940	43.204	ng	99	
89) Benzo(k)fluoranthene	15.139	252	378833	41.869	ng	99	
90) Benzo(a)pyrene	15.486	252	351174	40.866	ng	100	
91) Dibenzo(a,h)anthracene	17.115	278	302496	35.315	ng	98	
92) Benzo(g,h,i)perylene	17.568	276	274412	31.066	ng	97	

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

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