

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090823\
 Data File : BF135149.D
 Acq On : 08 Sep 2023 16:21
 Operator : CG\JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Sep 09 02:24:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 09 02:14:04 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							09/09/2023
1) 1,4-Dichlorobenzene-d4	6.763	152	106064	20.000	ng	0.00	Supervised By :mohammad
21) Naphthalene-d8	8.039	136	441982	20.000	ng	0.00	09/09/2023
39) Acenaphthene-d10	9.798	164	240863	20.000	ng	0.00	
64) Phenanthrene-d10	11.292	188	455545	20.000	ng	0.00	
76) Chrysene-d12	13.945	240	292177	20.000	ng	0.00	
86) Perylene-d12	15.404	264	211689	20.000	ng	0.00	09/10/2023
System Monitoring Compounds							
5) 2-Fluorophenol	5.387	112	75679	11.042	ng	0.00	
7) Phenol-d6	6.398	99	96383	11.176	ng	-0.01	
23) Nitrobenzene-d5	7.322	82	85801	10.567	ng	-0.01	
42) 2,4,6-Tribromophenol	10.586	330	30348	11.204	ng	-0.01	
45) 2-Fluorobiphenyl	9.122	172	179922	11.730	ng	0.00	
79) Terphenyl-d14	12.886	244	200631	10.573	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	2.475	88	15395	5.421	ng		96
3) Pyridine	3.246	79	38636	5.130	ng		94
4) n-Nitrosodimethylamine	3.175	42	18122	4.806	ng		93
6) Aniline	6.428	93	60282	5.691	ng		98
8) 2-Chlorophenol	6.545	128	40945	5.788	ng		97
9) Benzaldehyde	6.316	77	30481	6.303	ng		99
10) Phenol	6.410	94	51517	5.709	ng		85
11) bis(2-Chloroethyl)ether	6.498	93	36763	5.502	ng		99
12) 1,3-Dichlorobenzene	6.704	146	41737	5.774	ng		94
13) 1,4-Dichlorobenzene	6.781	146	42104	5.799	ng		98
14) 1,2-Dichlorobenzene	6.934	146	40068	5.895	ng		97
15) Benzyl Alcohol	6.904	79	34801	5.688	ng		97
16) 2,2'-oxybis(1-Chloropr...	7.039	45	62214	5.733	ng		99
17) 2-Methylphenol	7.016	107	35270	5.654	ng		96
18) Hexachloroethane	7.269	117	14862	5.584	ng		95
19) n-Nitroso-di-n-propyla...	7.169	70	29474	5.667	ng		98
20) 3+4-Methylphenols	7.169	107	44760	5.955	ng	#	75
22) Acetophenone	7.169	105	57229	5.614	ng	#	94
24) Nitrobenzene	7.339	77	44031	5.339	ng		97
25) Isophorone	7.581	82	79933	5.271	ng		99
26) 2-Nitrophenol	7.663	139	19888	4.941	ng		95
27) 2,4-Dimethylphenol	7.698	122	34687	5.455	ng		100
28) bis(2-Chloroethoxy)met...	7.792	93	48405	5.460	ng		98
29) 2,4-Dichlorophenol	7.904	162	34051	5.490	ng		93
30) 1,2,4-Trichlorobenzene	7.986	180	37118	5.488	ng		99
31) Naphthalene	8.063	128	123621	5.647	ng		99
32) Benzoic acid	7.769	122	20947m	4.271	ng		
33) 4-Chloroaniline	8.116	127	52738	5.530	ng		98
34) Hexachlorobutadiene	8.181	225	21301	5.473	ng		99
35) Caprolactam	8.451	113	10866	5.414	ng		96
36) 4-Chloro-3-methylphenol	8.598	107	37392	5.368	ng		99
37) 2-Methylnaphthalene	8.757	142	83472	5.620	ng		98
38) 1-Methylnaphthalene	8.851	142	78479	5.705	ng		97
40) 1,2,4,5-Tetrachloroben...	8.922	216	36915	5.403	ng		99
43) 2,4,6-Trichlorophenol	9.033	196	23822	5.254	ng		99
44) 2,4,5-Trichlorophenol	9.075	196	28589	5.450	ng		99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
46) 1,1'-Biphenyl	9.222	154	105170	5.628	ng	99	09/09/2023
47) 2-Chloronaphthalene	9.245	162	74457	5.367	ng	98	Supervised By :mohammad
48) 2-Nitroaniline	9.345	65	24765	5.053	ng	91	ahmed
49) Acenaphthylene	9.657	152	121620	5.672	ng	98	
50) Dimethylphthalate	9.522	163	93951	5.443	ng	100	
51) 2,6-Dinitrotoluene	9.586	165	19926	5.272	ng	96	
52) Acenaphthene	9.833	154	74247	5.517	ng	98	09/10/2023
53) 3-Nitroaniline	9.751	138	21900	5.143	ng	89	
55) Dibenzofuran	10.004	168	113402	5.800	ng	99	
56) 4-Nitrophenol	9.922	139	13857	4.900	ng	91	
57) 2,4-Dinitrotoluene	9.992	165	25366	5.356	ng	94	
58) Fluorene	10.345	166	91042	5.934	ng	99	
59) 2,3,4,6-Tetrachlorophenol	10.127	232	21693m	5.347	ng		
60) Diethylphthalate	10.222	149	90803	5.604	ng	99	
61) 4-Chlorophenyl-phenyle...	10.345	204	44611	5.913	ng	93	
62) 4-Nitroaniline	10.363	138	21445	5.152	ng	97	
63) Azobenzene	10.504	77	89497	5.563	ng	99	
65) 4,6-Dinitro-2-methylph...	10.404	198	10756	4.005	ng	92	
66) n-Nitrosodiphenylamine	10.463	169	78770	5.533	ng	96	
67) 4-Bromophenyl-phenylether	10.833	248	26196	5.413	ng	98	
68) Hexachlorobenzene	10.898	284	27780	5.357	ng	98	
69) Atrazine	10.992	200	23079	6.752	ng	96	
70) Pentachlorophenol	11.098	266	13247	4.707	ng	96	
71) Phenanthrene	11.316	178	136126	5.720	ng	99	
72) Anthracene	11.363	178	137456	5.658	ng	99	
73) Carbazole	11.521	167	115307	5.583	ng	98	
74) Di-n-butylphthalate	11.863	149	135392	5.502	ng	99	
75) Fluoranthene	12.510	202	136582	5.810	ng	96	
77) Benzidine	12.639	184	42322	7.147	ng	97	
78) Pyrene	12.739	202	140617	4.965	ng	99	
80) Butylbenzylphthalate	13.368	149	47993	4.850	ng	100	
81) Benzo(a)anthracene	13.933	228	108719	5.489	ng	99	
82) 3,3'-Dichlorobenzidine	13.898	252	31056	5.137	ng	# 93	
83) Chrysene	13.968	228	103295	5.341	ng	98	
84) Bis(2-ethylhexyl)phtha...	13.933	149	52326	5.009	ng	99	
85) Di-n-octyl phthalate	14.545	149	69254	4.319	ng	97	
87) Indeno(1,2,3-cd)pyrene	16.868	276	71457	4.646	ng	98	
88) Benzo(b)fluoranthene	14.980	252	71146	5.528	ng	98	
89) Benzo(k)fluoranthene	15.009	252	71261	5.702	ng	100	
90) Benzo(a)pyrene	15.339	252	63060	5.240	ng	98	
91) Dibenzo(a,h)anthracene	16.886	278	58199	4.636	ng	99	
92) Benzo(g,h,i)perylene	17.309	276	58624	4.526	ng	98	

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

