

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123021\
 Data File : BF126423.D
 Acq On : 30 Dec 2021 17:08
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 31 00:37:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 04:30:49 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh
 Patel

Supervised By :mohammad
 ahmed
 12/31/2021

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							12/31/2021
1) 1,4-Dichlorobenzene-d4	6.798	152	145857	20.000	ng	0.00	Supervised By :mohammad ahmed
21) Naphthalene-d8	8.081	136	590974	20.000	ng	0.00	
39) Acenaphthene-d10	9.839	164	271393	20.000	ng	0.00	
64) Phenanthrene-d10	11.322	188	425659	20.000	ng	0.00	
76) Chrysene-d12	13.963	240	226975	20.000	ng	# 0.00	
86) Perylene-d12	15.398	264	212035	20.000	ng	0.00	12/31/2021
System Monitoring Compounds							
5) 2-Fluorophenol	5.404	112	820433	80.777	ng	0.00	
7) Phenol-d6	6.434	99	1043363	79.312	ng	0.00	
23) Nitrobenzene-d5	7.363	82	919803	79.916	ng	0.00	
42) 2,4,6-Tribromophenol	10.627	330	173868	82.906	ng	0.00	
45) 2-Fluorobiphenyl	9.163	172	1105726	71.696	ng	0.00	
79) Terphenyl-d14	12.910	244	964258	82.542	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	2.493	88	219062	43.010	ng		99
3) Pyridine	3.222	79	591112	43.025	ng		98
4) n-Nitrosodimethylamine	3.169	42	223863	39.632	ng		99
6) Aniline	6.463	93	670838	41.010	ng		96
8) 2-Chlorophenol	6.587	128	412241	40.699	ng		93
9) Benzaldehyde	6.345	77	299320	36.755	ng		97
10) Phenol	6.445	94	597310	40.687	ng		99
11) bis(2-Chloroethyl)ether	6.534	93	452117	40.333	ng		99
12) 1,3-Dichlorobenzene	6.739	146	406401	40.151	ng		99
13) 1,4-Dichlorobenzene	6.816	146	410447	40.625	ng		98
14) 1,2-Dichlorobenzene	6.969	146	379668	38.567	ng		99
15) Benzyl Alcohol	6.939	79	376958	40.955	ng		99
16) 2,2'-oxybis(1-Chloropr...	7.075	45	703807	39.977	ng		99
17) 2-Methylphenol	7.051	107	359594	40.359	ng		99
18) Hexachloroethane	7.316	117	161491	40.118	ng		93
19) n-Nitroso-di-n-propyla...	7.216	70	287327	38.276	ng		98
20) 3+4-Methylphenols	7.204	107	403921	38.079	ng		99
22) Acetophenone	7.210	105	528973	38.718	ng		97
24) Nitrobenzene	7.381	77	459950	39.935	ng		99
25) Isophorone	7.622	82	814902	39.551	ng		99
26) 2-Nitrophenol	7.698	139	211147	41.355	ng		98
27) 2,4-Dimethylphenol	7.734	122	327532	41.026	ng		100
28) bis(2-Chloroethoxy)met...	7.833	93	532195	39.540	ng		98
29) 2,4-Dichlorophenol	7.939	162	294156	40.352	ng		99
30) 1,2,4-Trichlorobenzene	8.028	180	305702	40.237	ng		99
31) Naphthalene	8.104	128	1113813	39.930	ng		99
32) Benzoic acid	7.863	122	214646	38.674	ng		98
33) 4-Chloroaniline	8.151	127	478582	40.954	ng		99
34) Hexachlorobutadiene	8.228	225	146265	39.609	ng		99
35) Caprolactam	8.516	113	75519m	40.722	ng		
36) 4-Chloro-3-methylphenol	8.633	107	358351	40.280	ng		98
37) 2-Methylnaphthalene	8.798	142	656306	39.885	ng		98
38) 1-Methylnaphthalene	8.892	142	628805	39.525	ng		99
40) 1,2,4,5-Tetrachloroben...	8.963	216	235196	39.882	ng		99
41) Hexachlorocyclopentadiene	8.951	237	96997	40.398	ng		99
43) 2,4,6-Trichlorophenol	9.075	196	179023	41.048	ng		96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.110	196	192300	40.963	ng	99	12/31/2021
46) 1,1'-Biphenyl	9.263	154	775549	38.962	ng	99	Supervised By :mohammad ahmed
47) 2-Chloronaphthalene	9.286	162	593097	39.339	ng	99	
48) 2-Nitroaniline	9.380	65	244301	40.127	ng	89	
49) Acenaphthylene	9.698	152	914023	39.973	ng	99	
50) Dimethylphthalate	9.563	163	662558	39.106	ng	99	
51) 2,6-Dinitrotoluene	9.622	165	151294	41.349	ng	95	12/31/2021
52) Acenaphthene	9.875	154	588636	39.266	ng	98	
53) 3-Nitroaniline	9.786	138	201861	40.237	ng	99	
54) 2,4-Dinitrophenol	9.898	184	69459	39.095	ng	# 84	
55) Dibenzofuran	10.045	168	796245	39.290	ng	99	
56) 4-Nitrophenol	9.945	139	153243	42.951	ng	98	
57) 2,4-Dinitrotoluene	10.022	165	201376	42.659	ng	99	
58) Fluorene	10.386	166	553039	38.597	ng	99	
59) 2,3,4,6-Tetrachlorophenol	10.163	232	142259	42.220	ng	96	
60) Diethylphthalate	10.263	149	654000	40.053	ng	99	
61) 4-Chlorophenyl-phenyle...	10.380	204	245688	39.291	ng	93	
62) 4-Nitroaniline	10.404	138	192194	40.678	ng	93	
63) Azobenzene	10.539	77	825381	39.244	ng	99	
65) 4,6-Dinitro-2-methylph...	10.433	198	103164	44.755	ng	88	
66) n-Nitrosodiphenylamine	10.498	169	522510	39.566	ng	99	
67) 4-Bromophenyl-phenylether	10.869	248	155184	41.074	ng	99	
68) Hexachlorobenzene	10.933	284	179671	40.883	ng	98	
69) Atrazine	11.022	200	99770	43.874	ng	99	
70) Pentachlorophenol	11.127	266	93481	43.765	ng	99	
71) Phenanthrene	11.345	178	864643	39.732	ng	100	
72) Anthracene	11.398	178	870465	39.976	ng	100	
73) Carbazole	11.551	167	852024	39.959	ng	100	
74) Di-n-butylphthalate	11.886	149	964594	39.735	ng	99	
75) Fluoranthene	12.533	202	797870	40.171	ng	99	
77) Benzidine	12.651	184	176496	46.579	ng	100	
78) Pyrene	12.763	202	816703	42.578	ng	99	
80) Butylbenzylphthalate	13.380	149	358178	42.140	ng	99	
81) Benzo(a)anthracene	13.951	228	521461	39.949	ng	99	
82) 3,3'-Dichlorobenzidine	13.910	252	242986	39.901	ng	99	
83) Chrysene	13.986	228	541536	40.446	ng	100	
84) Bis(2-ethylhexyl)phtha...	13.945	149	368240	41.805	ng	98	
85) Di-n-octyl phthalate	14.557	149	666972	43.660	ng	99	
87) Indeno(1,2,3-cd)pyrene	16.821	276	596554	42.039	ng	100	
88) Benzo(b)fluoranthene	14.986	252	514323	40.718	ng	97	
89) Benzo(k)fluoranthene	15.009	252	476568	38.948	ng	99	
90) Benzo(a)pyrene	15.339	252	433007	40.884	ng	99	
91) Dibenzo(a,h)anthracene	16.839	278	491381	42.914	ng	99	
92) Benzo(g,h,i)perylene	17.251	276	510309	41.853	ng	98	

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

Manual Integrations

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