

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053124\
 Data File : BF138090.D
 Acq On : 31 May 2024 16:16
 Operator : CG\JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDICC050

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 06/01/2024

Supervised By :mohammad ahmed 06/01/2024

Quant Time: Jun 01 01:14:21 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF053124.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Jun 01 01:06:32 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.034	152	122528	20.000	ng	0.00
21) Naphthalene-d8	8.322	136	472080	20.000	ng	0.00
39) Acenaphthene-d10	10.086	164	265629	20.000	ng	0.00
64) Phenanthrene-d10	11.580	188	438052	20.000	ng	0.00
76) Chrysene-d12	14.239	240	252101	20.000	ng	0.00
86) Perylene-d12	15.804	264	249257	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.663	112	645945	94.241	ng	0.00
7) Phenol-d6	6.663	99	831035	94.154	ng	0.00
23) Nitrobenzene-d5	7.604	82	780413	96.329	ng	0.00
42) 2,4,6-Tribromophenol	10.880	330	291344	94.030	ng	0.00
45) 2-Fluorobiphenyl	9.398	172	1646189	91.864	ng	0.00
79) Terphenyl-d14	13.168	244	1681451	96.477	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.946	88	145622	48.936	ng	99
3) Pyridine	3.716	79	333820	50.752	ng	100
4) n-Nitrosodimethylamine	3.693	42	143377	49.642	ng	99
6) Aniline	6.710	93	198537m	48.007	ng	
8) 2-Chlorophenol	6.828	128	372149	47.329	ng	95
9) Benzaldehyde	6.592	77	148848	37.936	ng	99
10) Phenol	6.681	94	413789	47.266	ng	96
11) bis(2-Chloroethyl)ether	6.769	93	327963	45.344	ng	99
12) 1,3-Dichlorobenzene	6.981	146	435855	47.702	ng	99
13) 1,4-Dichlorobenzene	7.057	146	441707	48.090	ng	100
14) 1,2-Dichlorobenzene	7.210	146	410044	47.180	ng	99
15) Benzyl Alcohol	7.175	79	197315	49.427	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.304	45	382624	47.997	ng	99
17) 2-Methylphenol	7.287	107	298539	47.300	ng	98
18) Hexachloroethane	7.551	117	151299	48.492	ng	100
19) n-Nitroso-di-n-propyla...	7.451	70	237617	46.426	ng	98
20) 3+4-Methylphenols	7.439	107	373603	47.087	ng	# 84
22) Acetophenone	7.445	105	503442	47.335	ng	# 98
24) Nitrobenzene	7.622	77	390510	48.429	ng	98
25) Isophorone	7.857	82	666688	47.698	ng	99
26) 2-Nitrophenol	7.939	139	214750	49.828	ng	93
27) 2,4-Dimethylphenol	7.963	122	253650	49.167	ng	95
28) bis(2-Chloroethoxy)met...	8.063	93	406079	47.474	ng	99
29) 2,4-Dichlorophenol	8.181	162	335728	48.488	ng	99
30) 1,2,4-Trichlorobenzene	8.257	180	368540	48.029	ng	100
31) Naphthalene	8.345	128	1145936	47.178	ng	100
32) Benzoic acid	8.098	122	202431	49.607	ng	97
33) 4-Chloroaniline	8.457	127	395400m	48.289	ng	
34) Hexachlorobutadiene	8.451	225	235754	47.566	ng	98
35) Caprolactam	8.781	113	98863m	48.548	ng	
36) 4-Chloro-3-methylphenol	8.875	107	332677	47.928	ng	99
37) 2-Methylnaphthalene	9.039	142	753425	47.145	ng	99
38) 1-Methylnaphthalene	9.139	142	725352	46.640	ng	100
40) 1,2,4,5-Tetrachloroben...	9.204	216	362707	48.105	ng	100
41) Hexachlorocyclopentadiene	9.186	237	141375	53.751	ng	97
43) 2,4,6-Trichlorophenol	9.316	196	240706	48.127	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.363	196	265680	48.720	ng	98
46) 1,1'-Biphenyl	9.504	154	900066	47.111	ng	98
47) 2-Chloronaphthalene	9.533	162	722732	47.166	ng	99
48) 2-Nitroaniline	9.628	65	183308	47.774	ng	99
49) Acenaphthylene	9.951	152	1028492	46.901	ng	99
50) Dimethylphthalate	9.804	163	820689	46.625	ng	99
51) 2,6-Dinitrotoluene	9.869	165	191355	47.353	ng	94
52) Acenaphthene	10.122	154	645666	46.490	ng	99
53) 3-Nitroaniline	10.045	138	185083	48.118	ng	96
54) 2,4-Dinitrophenol	10.151	184	102657	52.251	ng #	88
55) Dibenzofuran	10.292	168	991637	46.503	ng	98
56) 4-Nitrophenol	10.204	139	132425	51.655	ng	95
57) 2,4-Dinitrotoluene	10.275	165	253450	47.562	ng	95
58) Fluorene	10.639	166	792832	46.147	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.410	232	206360	47.996	ng	98
60) Diethylphthalate	10.498	149	778805	46.584	ng	98
61) 4-Chlorophenyl-phenyle...	10.627	204	402380	46.381	ng	98
62) 4-Nitroaniline	10.663	138	150816	43.188	ng	98
63) Azobenzene	10.786	77	672094	46.783	ng	99
65) 4,6-Dinitro-2-methylph...	10.686	198	143351	52.188	ng	85
66) n-Nitrosodiphenylamine	10.745	169	665642	48.848	ng	100
67) 4-Bromophenyl-phenylether	11.122	248	251517	48.985	ng	98
68) Hexachlorobenzene	11.186	284	276689	48.496	ng	98
69) Atrazine	11.275	200	188360	50.067	ng	98
70) Pentachlorophenol	11.386	266	155544	52.661	ng	95
71) Phenanthrene	11.610	178	1048173	47.570	ng	99
72) Anthracene	11.663	178	1049476	47.757	ng	100
73) Carbazole	11.816	167	963002	47.166	ng	100
74) Di-n-butylphthalate	12.127	149	1160367	47.621	ng	100
75) Fluoranthene	12.804	202	1065515	45.831	ng	99
77) Benzidine	12.945	184	120624m	45.930	ng	
78) Pyrene	13.033	202	1064397	49.699	ng	99
80) Butylbenzylphthalate	13.639	149	377855	48.595	ng	99
81) Benzo(a)anthracene	14.227	228	789010	47.922	ng	98
82) 3,3'-Dichlorobenzidine	14.186	252	206949	46.922	ng	97
83) Chrysene	14.263	228	739277	47.555	ng	99
84) Bis(2-ethylhexyl)phtha...	14.192	149	506451	47.876	ng #	98
85) Di-n-octyl phthalate	14.815	149	760256	49.041	ng	99
87) Indeno(1,2,3-cd)pyrene	17.451	276	784452	52.214	ng	99
88) Benzo(b)fluoranthene	15.333	252	742132	48.330	ng	100
89) Benzo(k)fluoranthene	15.368	252	667557	45.754	ng	99
90) Benzo(a)pyrene	15.739	252	617963	48.143	ng	99
91) Dibenzo(a,h)anthracene	17.468	278	646138	52.422	ng	98
92) Benzo(g,h,i)perylene	17.951	276	662994	52.956	ng	99

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

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