

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011422\
 Data File : BP008813.D
 Acq On : 14 Jan 2022 17:47
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :

BNA_P

ClientSampleId :

ICVBP011422

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 01/17/2022
 Supervised By :Jagrut Upadhyay 01/17/2022

Quant Time: Jan 17 03:16:16 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jan 14 18:05:00 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	296228	20.000	ng	0.00
21) Naphthalene-d8	10.658	136	1124520	20.000	ng	0.00
39) Acenaphthene-d10	14.493	164	694712	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	1384903	20.000	ng	0.00
76) Chrysene-d12	21.339	240	1382425	20.000	ng	0.00
86) Perylene-d12	23.763	264	1478113	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1371674	71.606	ng	0.00
7) Phenol-d6	7.034	99	1666888	71.859	ng	0.00
23) Nitrobenzene-d5	9.016	82	1443609	72.883	ng	0.00
42) 2,4,6-Tribromophenol	15.987	330	751179	74.284	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	3798206	72.100	ng	0.00
79) Terphenyl-d14	19.810	244	5730135	69.965	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.346	88	273815	35.315	ng	98
3) Pyridine	3.746	79	595700	36.684	ng	97
4) n-Nitrosodimethylamine	3.652	42	276491	36.291	ng	# 85
6) Aniline	7.181	93	1035198	35.755	ng	98
8) 2-Chlorophenol	7.422	128	732639	35.911	ng	98
9) Benzaldehyde	6.993	77	555206	35.811	ng	99
10) Phenol	7.058	94	847885	35.853	ng	96
11) bis(2-Chloroethyl)ether	7.281	93	674891	35.869	ng	98
12) 1,3-Dichlorobenzene	7.746	146	861531	35.461	ng	99
13) 1,4-Dichlorobenzene	7.893	146	879569	35.721	ng	99
14) 1,2-Dichlorobenzene	8.211	146	837648	35.677	ng	99
15) Benzyl Alcohol	8.099	79	581253	35.718	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.387	45	760510	36.143	ng	98
17) 2-Methylphenol	8.305	107	567705	35.833	ng	97
18) Hexachloroethane	8.934	117	305675	36.576	ng	99
19) n-Nitroso-di-n-propyla...	8.669	70	489996	36.136	ng	96
20) 3+4-Methylphenols	8.634	107	753409	35.822	ng	99
22) Acetophenone	8.681	105	1033163	36.066	ng	# 97
24) Nitrobenzene	9.058	77	725043	36.239	ng	97
25) Isophorone	9.587	82	1298121	36.265	ng	98
26) 2-Nitrophenol	9.769	139	385523	36.653	ng	93
27) 2,4-Dimethylphenol	9.840	122	653077	36.416	ng	98
28) bis(2-Chloroethoxy)met...	10.069	93	849364	36.363	ng	99
29) 2,4-Dichlorophenol	10.310	162	705559	36.052	ng	100
30) 1,2,4-Trichlorobenzene	10.522	180	821970	35.856	ng	99
31) Naphthalene	10.710	128	2170619	35.539	ng	99
32) Benzoic acid	9.999	122	383036	37.216	ng	98
33) 4-Chloroaniline	10.816	127	894672	35.946	ng	98
34) Hexachlorobutadiene	10.999	225	513809	35.874	ng	99
35) Caprolactam	11.610	113	196156	36.359	ng	95
36) 4-Chloro-3-methylphenol	11.952	107	631751	35.782	ng	98
37) 2-Methylnaphthalene	12.322	142	1593961	35.699	ng	98
38) 1-Methylnaphthalene	12.540	142	1464018	35.723	ng	100
40) 1,2,4,5-Tetrachloroben...	12.693	216	910017	35.927	ng	99
41) Hexachlorocyclopentadiene	12.675	237	501425	38.687	ng	98
43) 2,4,6-Trichlorophenol	12.934	196	567245	36.220	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.010	196	616033	36.547	ng	98
46) 1,1'-Biphenyl	13.328	154	2026414	35.852	ng	99
47) 2-Chloronaphthalene	13.369	162	1567250	35.961	ng	98
48) 2-Nitroaniline	13.575	65	360075	37.210	ng	97
49) Acenaphthylene	14.216	152	2378270	36.119	ng	100
50) Dimethylphthalate	13.957	163	1879776	36.433	ng	99
51) 2,6-Dinitrotoluene	14.069	165	403359	36.861	ng	93
52) Acenaphthene	14.557	154	1415213	36.238	ng	100
53) 3-Nitroaniline	14.398	138	398863	37.063	ng	97
54) 2,4-Dinitrophenol	14.610	184	183776	41.527	ng #	90
55) Dibenzofuran	14.893	168	2310523	36.308	ng	98
56) 4-Nitrophenol	14.722	139	304941	39.350	ng	91
57) 2,4-Dinitrotoluene	14.863	165	542788	38.564	ng	93
58) Fluorene	15.545	166	1850664	36.820	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.128	232	510259m	36.749	ng	
60) Diethylphthalate	15.322	149	1839672	37.578	ng	98
61) 4-Chlorophenyl-phenyle...	15.540	204	1007283	36.788	ng	99
62) 4-Nitroaniline	15.563	138	404047	38.766	ng	90
63) Azobenzene	15.834	77	1535695	37.784	ng	96
65) 4,6-Dinitro-2-methylph...	15.628	198	279566	40.405	ng	90
66) n-Nitrosodiphenylamine	15.751	169	1597613	35.843	ng	99
67) 4-Bromophenyl-phenylether	16.434	248	625256	36.156	ng	97
68) Hexachlorobenzene	16.557	284	701219	35.386	ng	98
69) Atrazine	16.716	200	617279	37.267	ng	98
70) Pentachlorophenol	16.904	266	409263	38.772	ng	100
71) Phenanthrene	17.292	178	2800558	35.916	ng	99
72) Anthracene	17.381	178	2878501	36.799	ng	99
73) Carbazole	17.657	167	2519123	37.454	ng	99
74) Di-n-butylphthalate	18.210	149	3166933	39.126	ng	99
75) Fluoranthene	19.263	202	3291044	37.954	ng	96
77) Benzidine	19.439	184	1974038	40.420	ng	98
78) Pyrene	19.616	202	3339171	34.617	ng	98
80) Butylbenzylphthalate	20.486	149	1425029	36.733	ng	99
81) Benzo(a)anthracene	21.328	228	3309189	35.581	ng	98
82) 3,3'-Dichlorobenzidine	21.257	252	1467920	36.534	ng	99
83) Chrysene	21.380	228	3218025	35.694	ng	97
84) Bis(2-ethylhexyl)phtha...	21.251	149	2231433	37.142	ng	97
85) Di-n-octyl phthalate	22.180	149	3794372	37.335	ng	99
87) Indeno(1,2,3-cd)pyrene	26.292	276	4327481	35.530	ng	98
88) Benzo(b)fluoranthene	23.022	252	3496914	35.425	ng	99
89) Benzo(k)fluoranthene	23.075	252	3490800	36.533	ng	98
90) Benzo(a)pyrene	23.657	252	3450558	36.215	ng	99
91) Dibenzo(a,h)anthracene	26.310	278	3677633	35.509	ng	99
92) Benzo(g,h,i)perylene	27.068	276	3534991	34.958	ng	98

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

