

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080322\
 Data File : BF129546.D
 Acq On : 03 Aug 2022 18:17
 Operator : CG\JU
 Sample : N3930-05MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-98MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/04/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

Quant Time: Aug 04 02:19:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	140308	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	521066	20.000	ng	-0.01
39) Acenaphthene-d10	9.910	164	225139	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	287610	20.000	ng	0.00
76) Chrysene-d12	14.045	240	262225	20.000	ng	0.00
86) Perylene-d12	15.527	264	324886	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	782204	90.815	ng	0.00
7) Phenol-d6	6.516	99	945521	87.336	ng	0.00
23) Nitrobenzene-d5	7.434	82	600934	62.956	ng	-0.01
42) 2,4,6-Tribromophenol	10.704	330	160915	74.341	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	996944	68.501	ng	0.00
79) Terphenyl-d14	12.980	244	668061	48.064	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.699	88	176091	45.313	ng	89
3) Pyridine	3.481	79	450337	41.729	ng	89
4) n-Nitrosodimethylamine	3.434	42	230312	48.599	ng	# 94
6) Aniline	6.534	93	356298	26.473	ng	# 43
8) 2-Chlorophenol	6.663	128	464522	49.365	ng	95
9) Benzaldehyde	6.422	77	217005	29.516	ng	98
10) Phenol	6.528	94	542116m	47.022	ng	
11) bis(2-Chloroethyl)ether	6.604	93	440376	48.165	ng	92
12) 1,3-Dichlorobenzene	6.810	146	495667	49.114	ng	99
13) 1,4-Dichlorobenzene	6.887	146	503501	49.056	ng	99
14) 1,2-Dichlorobenzene	7.040	146	471698	49.998	ng	98
15) Benzyl Alcohol	7.016	79	382121	48.667	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.145	45	607367	44.192	ng	80
17) 2-Methylphenol	7.134	107	347110	44.464	ng	# 93
18) Hexachloroethane	7.381	117	188145	48.682	ng	92
19) n-Nitroso-di-n-propyla...	7.281	70	302909	46.217	ng	94
20) 3+4-Methylphenols	7.287	107	433270	43.651	ng	99
22) Acetophenone	7.281	105	612489	50.488	ng	# 97
24) Nitrobenzene	7.457	77	471278	47.946	ng	94
25) Isophorone	7.692	82	810698	47.381	ng	98
26) 2-Nitrophenol	7.769	139	234456	48.350	ng	99
27) 2,4-Dimethylphenol	7.810	122	386212m	53.097	ng	
28) bis(2-Chloroethoxy)met...	7.898	93	506381	51.018	ng	98
29) 2,4-Dichlorophenol	8.016	162	341107	45.435	ng	100
30) 1,2,4-Trichlorobenzene	8.092	180	362586	46.660	ng	98
31) Naphthalene	8.175	128	1241071	46.880	ng	99
32) Benzoic acid	7.934	122	171398	32.595	ng	100
33) 4-Chloroaniline	8.228	127	176706	16.258	ng	97
34) Hexachlorobutadiene	8.287	225	218905	49.557	ng	98
35) Caprolactam	8.598	113	89714m	36.756	ng	
36) 4-Chloro-3-methylphenol	8.716	107	333878	41.931	ng	97
37) 2-Methylnaphthalene	8.869	142	772490	44.902	ng	100
38) 1-Methylnaphthalene	8.963	142	727049	43.778	ng	98
40) 1,2,4,5-Tetrachloroben...	9.034	216	325996	55.144	ng	97
41) Hexachlorocyclopentadiene	9.016	237	301517	114.540	ng	99
43) 2,4,6-Trichlorophenol	9.151	196	206147	48.011	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	212227	45.451	ng	94
46) 1,1'-Biphenyl	9.328	154	890019	54.168	ng	99
47) 2-Chloronaphthalene	9.357	162	690977	52.021	ng	98
48) 2-Nitroaniline	9.457	65	200536	45.404	ng	94
49) Acenaphthylene	9.775	152	992319	49.167	ng	99
50) Dimethylphthalate	9.628	163	729321	46.926	ng	99
51) 2,6-Dinitrotoluene	9.698	165	160342	46.094	ng	92
52) Acenaphthene	9.945	154	666991m	50.598	ng	
53) 3-Nitroaniline	9.869	138	113726	27.686	ng	# 89
54) 2,4-Dinitrophenol	9.981	184	126830	71.420	ng	# 88
55) Dibenzofuran	10.116	168	862058	47.439	ng	97
56) 4-Nitrophenol	10.045	139	187647	61.921	ng	96
57) 2,4-Dinitrotoluene	10.104	165	197337	43.425	ng	88
58) Fluorene	10.457	166	654393	45.007	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	139947	38.784	ng	95
60) Diethylphthalate	10.328	149	642884	42.790	ng	100
61) 4-Chlorophenyl-phenyle...	10.451	204	298815	46.617	ng	97
62) 4-Nitroaniline	10.480	138	132979	32.634	ng	98
63) Azobenzene	10.610	77	631043	41.179	ng	95
65) 4,6-Dinitro-2-methylph...	10.510	198	82347	45.362	ng	92
66) n-Nitrosodiphenylamine	10.569	169	527773	55.102	ng	96
67) 4-Bromophenyl-phenylether	10.939	248	170477	54.130	ng	95
68) Hexachlorobenzene	11.010	284	183151	53.671	ng	93
69) Atrazine	11.092	200	138356	47.557	ng	95
70) Pentachlorophenol	11.210	266	134500	72.501	ng	99
71) Phenanthrene	11.422	178	766148	48.800	ng	98
72) Anthracene	11.475	178	764502	49.552	ng	99
73) Carbazole	11.633	167	670249	46.149	ng	98
74) Di-n-butylphthalate	11.951	149	751304	43.439	ng	99
75) Fluoranthene	12.616	202	736370	46.291	ng	99
77) Benzidine	12.733	184	173521	18.731	ng	97
78) Pyrene	12.845	202	764133	34.847	ng	97
80) Butylbenzylphthalate	13.457	149	341377	35.892	ng	90
81) Benzo(a)anthracene	14.033	228	826999	46.169	ng	100
82) 3,3'-Dichlorobenzidine	13.998	252	212831	35.987	ng	99
83) Chrysene	14.074	228	792444	46.941	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	497100	39.699	ng	99
85) Di-n-octyl phthalate	14.621	149	958025	53.167	ng	96
87) Indeno(1,2,3-cd)pyrene	17.045	276	1047698	47.888	ng	100
88) Benzo(b)fluoranthene	15.092	252	1012026	46.960	ng	99
89) Benzo(k)fluoranthene	15.127	252	907901	42.990	ng	99
90) Benzo(a)pyrene	15.468	252	879976	50.301	ng	98
91) Dibenzo(a,h)anthracene	17.057	278	897297	50.391	ng	99
92) Benzo(g,h,i)perylene	17.504	276	863458	47.454	ng	99

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

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