

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021519\
 Data File : BM018823.D
 Acq On : 15 Feb 2019 18:55
 Operator : JU/SJ
 Sample : K1501-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CH-021419MSD

Manual Integrations
 APPROVED

Sohil
 2/18/2019 12:09:26 PM

Quant Time: Feb 15 23:19:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	331983	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1354592	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	765230	20.00	ng	0.00
64) Phenanthrene-d10	17.11	188	1682792	20.00	ng	0.00
76) Chrysene-d12	21.31	240	1680386	20.00	ng	0.00
87) Perylene-d12	23.56	264	1417024	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	4197176	207.15	ng	0.00
7) Phenol-d6	6.89	99	5976129	209.37	ng	0.00
23) Nitrobenzene-d5	8.88	82	3846616	131.30	ng	0.00
42) 2,4,6-Tribromophenol	15.86	330	2327932	211.04	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	7147506	123.99	ng	0.00
79) Terphenyl-d14	19.74	244	10600522	130.14	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.25	88	413436	46.849	ng	98
3) Pyridine	3.65	79	1236475	51.366	ng	100
4) n-Nitrosodimethylamine	3.57	42	426655	69.672	ng	99
6) Aniline	7.06	93	1680629	48.055	ng	99
8) 2-Chlorophenol	7.28	128	1418883	63.783	ng	99
9) Benzaldehyde	6.86	77	821028	59.940	ng	98
10) Phenol	6.92	94	2082459	69.338	ng	99
11) bis(2-Chloroethyl)ether	7.15	93	1361625	59.433	ng	99
12) 1,3-Dichlorobenzene	7.59	146	1411059	56.411	ng	99
13) 1,4-Dichlorobenzene	7.75	146	1448008	56.862	ng	100
14) 1,2-Dichlorobenzene	8.06	146	1398258	57.163	ng	99
15) Benzyl Alcohol	7.96	79	1433790	63.217	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.24	45	1305150	58.860	ng	99
17) 2-Methylphenol	8.16	107	1235588	64.638	ng	99
18) Hexachloroethane	8.77	117	544171	58.538	ng	97
19) n-Nitroso-di-n-propylamine	8.52	70	1309129	59.235	ng	99
20) 3+4-Methylphenols	8.49	107	1691323	64.075	ng	99
22) Acetophenone	8.54	105	2159696	58.057	ng	# 99
24) Nitrobenzene	8.92	77	1855902	61.031	ng	100
25) Isophorone	9.44	82	3265029	69.849	ng	100
26) 2-Nitrophenol	9.62	139	776045	64.083	ng	98
27) 2,4-Dimethylphenol	9.68	122	1300916	73.551	ng	99
28) bis(2-Chloroethoxy)methane	9.92	93	1911411	63.234	ng	99
29) 2,4-Dichlorophenol	10.15	162	1343707	66.162	ng	99
30) 1,2,4-Trichlorobenzene	10.36	180	1379499	58.816	ng	98
31) Naphthalene	10.55	128	4233822	64.086	ng	99
32) Benzoic acid	9.86	122	859007	55.676	ng	97
33) 4-Chloroaniline	10.67	127	445007	15.069	ng	98
34) Hexachlorobutadiene	10.81	225	754207	55.400	ng	99
35) Caprolactam	11.51	113	451399m	54.461	ng	
36) 4-Chloro-3-methylphenol	11.80	107	1566742	63.985	ng	98
37) 2-Methylnaphthalene	12.16	142	3124180	67.167	ng	99
38) 1-Methylnaphthalene	12.39	142	2974860	65.404	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.53	216	1480290	60.472	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021519\
 Data File : BM018823.D
 Acq On : 15 Feb 2019 18:55
 Operator : JU/SJ
 Sample : K1501-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CH-021419MSD

Manual Integrations
 APPROVED

Sohil
 2/18/2019 12:09:26 PM

Quant Time: Feb 15 23:19:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.50	237	1704395	136.184	nq	99
43) 2,4,6-Trichlorophenol	12.79	196	1080106	66.652	nq	98
44) 2,4,5-Trichlorophenol	12.86	196	1145663	67.451	nq	99
46) 1,1'-Biphenyl	13.19	154	3945188	59.894	nq	100
47) 2-Chloronaphthalene	13.23	162	3080809	60.555	nq	99
48) 2-Nitroaniline	13.46	65	1103882	65.330	nq	99
49) Acenaphthylene	14.08	152	5021101	66.280	nq	99
50) Dimethylphthalate	13.83	163	4390875	68.850	nq	99
51) 2,6-Dinitrotoluene	13.96	165	883696	63.967	nq	97
52) Acenaphthene	14.43	154	2905276	62.544	nq	100
53) 3-Nitroaniline	14.29	138	446563	28.609	nq	99
54) 2,4-Dinitrophenol	14.50	184	1033238	115.142	nq	98
55) Dibenzofuran	14.76	168	4691020	61.441	nq	99
56) 4-Nitrophenol	14.61	139	1556008	124.874	nq	97
57) 2,4-Dinitrotoluene	14.75	165	1260205	66.207	nq	98
58) Fluorene	15.41	166	3861774	66.114	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.99	232	1026074	68.344	nq	100
60) Diethylphthalate	15.19	149	4047982	61.533	nq	99
61) 4-Chlorophenyl-phenylether	15.41	204	1934759	59.077	nq	96
62) 4-Nitroaniline	15.46	138	897917	58.677	nq	98
63) Azobenzene	15.70	77	4369368	61.945	nq	98
65) 4,6-Dinitro-2-methylphenol	15.50	198	706656	59.810	nq	99
66) n-Nitrosodiphenylamine	15.63	169	3353206	63.070	nq	99
67) 4-Bromophenyl-phenylether	16.30	248	1150502	61.080	nq	99
68) Hexachlorobenzene	16.40	284	1330083	60.751	nq	99
69) Atrazine	16.59	200	1112232	64.891	nq	100
70) Pentachlorophenol	16.76	266	1776196	125.998	nq	100
71) Phenanthrene	17.16	178	5920664	65.468	nq	100
72) Anthracene	17.24	178	6071120	68.971	nq	99
73) Carbazole	17.52	167	5556308	60.418	nq	100
74) Di-n-butylphthalate	18.07	149	6992468	66.113	nq	99
75) Fluoranthene	19.17	202	6815852	65.245	nq	97
77) Benzidine	19.37	184	3836387	101.372	nq	100
78) Pyrene	19.54	202	6946715	71.179	nq	99
80) Butylbenzylphthalate	20.44	149	3326418	74.788	nq	97
81) Benzo(a)anthracene	21.29	228	6572153	67.095	nq	99
82) 3,3'-Dichlorobenzidine	21.23	252	1444996	37.326	nq	99
83) Chrysene	21.34	228	6150488	65.510	nq	100
84) Bis(2-ethylhexyl)phthalate	21.21	149	4783671	73.484	nq	100
85) Di-n-octyl phthalate	22.10	149	7937460	72.547	nq	100
86) Indeno(1,2,3-cd)pyrene	25.86	276	5807635	53.578	nq	99
88) Benzo(b)fluoranthene	22.89	252	5783735	71.340	nq	100
89) Benzo(k)fluoranthene	22.93	252	5862868	72.638	nq	100
90) Benzo(a)pyrene	23.47	252	5391793	70.203	nq	99
91) Dibenzo(a,h)anthracene	25.87	278	4960935	65.197	nq	99
92) Benzo(g,h,i)perylene	26.56	276	4616041	65.162	nq	100

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

