

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021219\
 Data File : BM018801.D
 Acq On : 13 Feb 2019 11:57
 Operator : JU/SJ
 Sample : PB117038BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB117038BSD

Manual Integrations
 APPROVED

Sohil
 2/13/2019 1:37:45 PM

Quant Time: Feb 13 13:03:18 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	288697	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1242784	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	741715	20.00	ng	0.00
64) Phenanthrene-d10	17.12	188	1697411	20.00	ng	0.00
76) Chrysene-d12	21.31	240	1847288	20.00	ng	0.00
87) Perylene-d12	23.56	264	1752659	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	1403692	79.67	ng	0.00
7) Phenol-d6	6.89	99	2061867	83.07	ng	0.00
23) Nitrobenzene-d5	8.88	82	2198494	81.80	ng	0.00
42) 2,4,6-Tribromophenol	15.86	330	885554	82.83	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	4459421	79.81	ng	0.00
79) Terphenyl-d14	19.75	244	8157780	91.10	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.25	88	315608	41.125	ng	99
3) Pyridine	3.65	79	966547	46.173	ng	99
4) n-Nitrosodimethylamine	3.57	42	267340	50.202	ng	95
6) Aniline	7.05	93	1420382	46.703	ng	99
8) 2-Chlorophenol	7.28	128	845635	43.713	ng	98
9) Benzaldehyde	6.87	77	625393	48.847	ng	99
10) Phenol	6.91	94	1162548	44.512	ng	99
11) bis(2-Chloroethyl)ether	7.15	93	857104	43.021	ng	100
12) 1,3-Dichlorobenzene	7.59	146	900263	41.386	ng	99
13) 1,4-Dichlorobenzene	7.75	146	919199	41.508	ng	99
14) 1,2-Dichlorobenzene	8.06	146	888272	41.759	ng	99
15) Benzyl Alcohol	7.96	79	898009	45.531	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.24	45	815716	42.303	ng	98
17) 2-Methylphenol	8.15	107	747097	44.943	ng	99
18) Hexachloroethane	8.78	117	341469	42.240	ng	98
19) n-Nitroso-di-n-propylamine	8.52	70	856360	44.558	ng	100
20) 3+4-Methylphenols	8.49	107	1029010	44.829	ng	98
22) Acetophenone	8.54	105	1420659	41.626	ng	# 99
24) Nitrobenzene	8.92	77	1188885	42.614	ng	100
25) Isophorone	9.44	82	2152715	50.196	ng	99
26) 2-Nitrophenol	9.62	139	481202	43.311	ng	99
27) 2,4-Dimethylphenol	9.68	122	808044	49.795	ng	99
28) bis(2-Chloroethoxy)methane	9.92	93	1234879	44.528	ng	100
29) 2,4-Dichlorophenol	10.15	162	824952	44.274	ng	99
30) 1,2,4-Trichlorobenzene	10.36	180	879289	40.862	ng	98
31) Naphthalene	10.55	128	2721098	44.894	ng	99
32) Benzoic acid	9.84	122	570324m	40.290	ng	
33) 4-Chloroaniline	10.68	127	1222139	45.107	ng	99
34) Hexachlorobutadiene	10.82	225	483576	38.717	ng	99
35) Caprolactam	11.50	113	301045m	39.588	ng	
36) 4-Chloro-3-methylphenol	11.80	107	998183	44.432	ng	99
37) 2-Methylnaphthalene	12.17	142	2035952	47.709	ng	99
38) 1-Methylnaphthalene	12.39	142	1933819	46.341	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.53	216	966259	40.724	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.50	237	1243786	102.531	nq	99
43) 2,4,6-Trichlorophenol	12.79	196	688449	43.830	nq	99
44) 2,4,5-Trichlorophenol	12.86	196	725465	44.066	nq	99
46) 1,1'-Biphenyl	13.19	154	2585329	40.494	nq	100
47) 2-Chloronaphthalene	13.23	162	2016071	40.883	nq	99
48) 2-Nitroaniline	13.46	65	728751	44.496	nq	98
49) Acenaphthylene	14.08	152	3362072	45.787	nq	99
50) Dimethylphthalate	13.83	163	2645360	42.795	nq	99
51) 2,6-Dinitrotoluene	13.96	165	587866	43.902	nq	97
52) Acenaphthene	14.43	154	1917052	42.578	nq	99
53) 3-Nitroaniline	14.29	138	629030	41.576	nq	99
54) 2,4-Dinitrophenol	14.50	184	509599	61.419	nq	97
55) Dibenzofuran	14.76	168	3127332	42.259	nq	100
56) 4-Nitrophenol	14.60	139	1031165	85.378	nq	99
57) 2,4-Dinitrotoluene	14.75	165	837155	45.376	nq	99
58) Fluorene	15.41	166	2566058	45.324	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.99	232	678382	46.618	nq	99
60) Diethylphthalate	15.19	149	2733863	42.875	nq	99
61) 4-Chlorophenyl-phenylether	15.41	204	1272448	40.086	nq	98
62) 4-Nitroaniline	15.46	138	634922	42.806	nq	100
63) Azobenzene	15.70	77	2933914	42.913	nq	98
65) 4,6-Dinitro-2-methylphenol	15.50	198	405472	34.023	nq	99
66) n-Nitrosodiphenylamine	15.63	169	2245267	41.867	nq	100
67) 4-Bromophenyl-phenylether	16.30	248	766822	40.360	nq	98
68) Hexachlorobenzene	16.41	284	896311	40.586	nq	95
69) Atrazine	16.59	200	774702	44.809	nq	99
70) Pentachlorophenol	16.76	266	1134026	79.751	nq	99
71) Phenanthrene	17.16	178	4076849	44.692	nq	100
72) Anthracene	17.25	178	4187591	47.164	nq	100
73) Carbazole	17.53	167	3850569	41.510	nq	99
74) Di-n-butylphthalate	18.08	149	4745452	44.482	nq	100
75) Fluoranthene	19.17	202	4788734	45.446	nq	98
77) Benzidine	19.38	184	5429129	130.497	nq	100
78) Pyrene	19.54	202	4919727	45.855	nq	100
80) Butylbenzylphthalate	20.45	149	2272242	46.471	nq	100
81) Benzo(a)anthracene	21.29	228	4892203	45.432	nq	100
82) 3,3'-Dichlorobenzidine	21.23	252	1877007	44.105	nq	99
83) Chrysene	21.34	228	4575273	44.329	nq	100
84) Bis(2-ethylhexyl)phthalate	21.21	149	3304240	46.172	nq	99
85) Di-n-octyl phthalate	22.10	149	5651326	46.985	nq	100
86) Indeno(1,2,3-cd)pyrene	25.86	276	4705516	39.488	nq	100
88) Benzo(b)fluoranthene	22.89	252	4817244	48.040	nq	100
89) Benzo(k)fluoranthene	22.93	252	4591711	45.995	nq	100
90) Benzo(a)pyrene	23.47	252	4442381	46.765	nq	98
91) Dibenzo(a,h)anthracene	25.87	278	4031191	42.833	nq	99
92) Benzo(g,h,i)perylene	26.56	276	3665943	41.840	nq	100

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

