

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092519\
 Data File : BM022781.D
 Acq On : 25 Sep 2019 20:57
 Operator : JU
 Sample : K4980-04MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SAND-STOCKPILE-1MSD

Manual Integrations
 APPROVED

mohammad
 9/26/2019 3:55:10 PM

Quant Time: Sep 26 06:48:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM092019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 16:34:01 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	242993	20.00	ng	0.00
21) Naphthalene-d8	10.61	136	873983	20.00	ng	-0.01
39) Acenaphthene-d10	14.46	164	478058	20.00	ng	0.00
64) Phenanthrene-d10	17.19	188	893833	20.00	ng	0.00
76) Chrysene-d12	21.37	240	926533	20.00	ng	0.00
87) Perylene-d12	23.64	264	1099903	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1432319	93.48	ng	0.00
7) Phenol-d6	6.99	99	1875507	95.31	ng	0.00
23) Nitrobenzene-d5	8.97	82	1067399	67.79	ng	0.00
42) 2,4,6-Tribromophenol	15.94	330	637215	106.78	ng	0.00
45) 2-Fluorobiphenyl	13.08	172	2431246	70.03	ng	0.00
79) Terphenyl-d14	19.82	244	3136726	65.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.31	88	166809	27.057	ng	# 100
3) Pyridine	3.72	79	400891	24.670	ng	100
4) n-Nitrosodimethylamine	3.62	42	193475	32.705	ng	# 99
6) Aniline	7.15	93	608253	25.778	ng	100
8) 2-Chlorophenol	7.39	128	555166	35.564	ng	100
9) Benzaldehyde	6.96	77	352507	31.460	ng	100
10) Phenol	7.02	94	660841	35.670	ng	99
11) bis(2-Chloroethyl)ether	7.25	93	499009	34.283	ng	99
12) 1,3-Dichlorobenzene	7.71	146	606576	32.511	ng	100
13) 1,4-Dichlorobenzene	7.86	146	624954	33.087	ng	99
14) 1,2-Dichlorobenzene	8.17	146	586469	32.568	ng	99
15) Benzyl Alcohol	8.06	79	428318	34.784	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.35	45	663857	31.397	ng	99
17) 2-Methylphenol	8.26	107	425897	34.731	ng	99
18) Hexachloroethane	8.90	117	220278	31.797	ng	97
19) n-Nitroso-di-n-propylamine	8.63	70	367562	33.448	ng	98
20) 3+4-Methylphenols	8.59	107	563694	34.512	ng	99
22) Acetophenone	8.64	105	738486	34.231	ng	99
24) Nitrobenzene	9.02	77	542569	36.017	ng	98
25) Isophorone	9.55	82	981594	37.917	ng	99
26) 2-Nitrophenol	9.73	139	275392	42.347	ng	97
27) 2,4-Dimethylphenol	9.79	122	473227	43.467	ng	100
28) bis(2-Chloroethoxy)methane	10.02	93	646654	36.097	ng	100
29) 2,4-Dichlorophenol	10.26	162	485093	39.410	ng	98
30) 1,2,4-Trichlorobenzene	10.47	180	543520	35.890	ng	99
31) Naphthalene	10.66	128	1633617	37.829	ng	100
32) Benzoic acid	9.90	122	226304	31.710	ng	95
33) 4-Chloroaniline	10.77	127	341822	18.231	ng	99
34) Hexachlorobutadiene	10.96	225	332461	36.828	ng	99
35) Caprolactam	11.55	113	132028m	36.181	ng	
36) 4-Chloro-3-methylphenol	11.89	107	454995	37.888	ng	99
37) 2-Methylnaphthalene	12.27	142	1154402	38.404	ng	100
38) 1-Methylnaphthalene	12.49	142	1071370	37.548	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.65	216	583815	36.041	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.63	237	402675	45.561	nq	99
43) 2,4,6-Trichlorophenol	12.89	196	372045	39.794	nq	99
44) 2,4,5-Trichlorophenol	12.96	196	388445	39.871	nq	98
46) 1,1'-Biphenyl	13.29	154	1396428	35.294	nq	100
47) 2-Chloronaphthalene	13.33	162	1071778	36.344	nq	99
48) 2-Nitroaniline	13.53	65	269135	35.048	nq	99
49) Acenaphthylene	14.17	152	1800856	41.238	nq	100
50) Dimethylphthalate	13.91	163	1534409	44.414	nq	100
51) 2,6-Dinitrotoluene	14.02	165	273820	37.891	nq	99
52) Acenaphthene	14.52	154	990823	34.680	nq	100
53) 3-Nitroaniline	14.35	138	205080	24.651	nq	99
54) 2,4-Dinitrophenol	14.56	184	153368	42.717	nq	96
55) Dibenzofuran	14.85	168	1535624	36.817	nq	99
56) 4-Nitrophenol	14.66	139	481264	75.263	nq	97
57) 2,4-Dinitrotoluene	14.81	165	355592	37.560	nq	99
58) Fluorene	15.50	166	1262711	37.255	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.08	232	328662	38.681	nq	97
60) Diethylphthalate	15.27	149	1246511	37.107	nq	100
61) 4-Chlorophenyl-phenylether	15.50	204	635864	36.512	nq	99
62) 4-Nitroaniline	15.51	138	283795	32.537	nq	98
63) Azobenzene	15.79	77	1084622	35.754	nq	99
65) 4,6-Dinitro-2-methylphenol	15.57	198	115829	28.864	nq	96
66) n-Nitrosodiphenylamine	15.71	169	1072704	39.284	nq	100
67) 4-Bromophenyl-phenylether	16.39	248	388011	39.029	nq	99
68) Hexachlorobenzene	16.51	284	421064	36.318	nq	96
69) Atrazine	16.66	200	320670	36.124	nq	99
70) Pentachlorophenol	16.84	266	513427	83.039	nq	99
71) Phenanthrene	17.23	178	2071048	41.064	nq	100
72) Anthracene	17.33	178	1958713	40.072	nq	99
73) Carbazole	17.59	167	1676614	36.861	nq	99
74) Di-n-butylphthalate	18.16	149	2106781	40.849	nq	100
75) Fluoranthene	19.25	202	2422730	42.569	nq	98
77) Benzidine	19.43	184	1248194	49.107	nq	100
78) Pyrene	19.61	202	2716572	44.739	nq	99
80) Butylbenzylphthalate	20.51	149	934153	40.480	nq	99
81) Benzo(a)anthracene	21.36	228	2512184	41.835	nq	98
82) 3,3'-Dichlorobenzidine	21.29	252	824394	38.712	nq	100
83) Chrysene	21.41	228	2414324	41.471	nq	100
84) Bis(2-ethylhexyl)phthalate	21.29	149	1438363	42.067	nq	99
85) Di-n-octyl phthalate	22.18	149	2474917	45.415	nq	98
86) Indeno(1,2,3-cd)pyrene	25.96	276	3012967	43.361	nq	# 100
88) Benzo(b)fluoranthene	22.96	252	2706398	40.639	nq	100
89) Benzo(k)fluoranthene	23.01	252	2409949	36.471	nq	99
90) Benzo(a)pyrene	23.55	252	2596267	42.308	nq	99
91) Dibenzo(a,h)anthracene	25.97	278	2412892	37.591	nq	98
92) Benzo(g,h,i)perylene	26.67	276	2529190	41.485	nq	98

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

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