

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000881.D
 Acq On : 18 Oct 2019 17:26
 Operator : JU
 Sample : K5401-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR200031-RT4221MSD

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:39:15 AM

Quant Time: Oct 19 04:32:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 16:40:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	151912	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	545216	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	306422	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	574781	20.00	ng	0.00
76) Chrysene-d12	13.96	240	483886	20.00	ng	0.00
87) Perylene-d12	15.44	264	552544	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	940644	99.53	ng	0.00
7) Phenol-d6	6.39	99	1187661	104.29	ng	0.00
23) Nitrobenzene-d5	7.30	82	665115	74.51	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	396880	121.55	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1438255	75.94	ng	0.00
79) Terphenyl-d14	12.89	244	1676063	70.13	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.45	88	103238	27.356	ng	99
3) Pyridine	3.19	79	261606	25.371	ng	99
4) n-Nitrosodimethylamine	3.12	42	90260	31.108	ng	100
6) Aniline	6.40	93	264767	19.137	ng	# 31
8) 2-Chlorophenol	6.53	128	331178	35.508	ng	98
9) Benzaldehyde	6.29	77	166104	28.001	ng	98
10) Phenol	6.40	94	405428	34.770	ng	87
11) bis(2-Chloroethyl)ether	6.48	93	312383	34.822	ng	100
12) 1,3-Dichlorobenzene	6.68	146	371965	32.899	ng	99
13) 1,4-Dichlorobenzene	6.76	146	382314	33.605	ng	99
14) 1,2-Dichlorobenzene	6.91	146	353996	33.702	ng	99
15) Benzyl Alcohol	6.89	79	258011	35.081	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.02	45	262502	31.723	ng	99
17) 2-Methylphenol	7.00	107	269917	35.017	ng	98
18) Hexachloroethane	7.25	117	129574	32.001	ng	98
19) n-Nitroso-di-n-propylamine	7.16	70	187823	35.781	ng	97
20) 3+4-Methylphenols	7.16	107	347844	38.566	ng	91
22) Acetophenone	7.15	105	456356	36.224	ng	# 97
24) Nitrobenzene	7.32	77	305536	35.487	ng	98
25) Isophorone	7.56	82	567845	35.841	ng	99
26) 2-Nitrophenol	7.64	139	175494	38.760	ng	98
27) 2,4-Dimethylphenol	7.69	122	283595	40.861	ng	100
28) bis(2-Chloroethoxy)methane	7.77	93	376461	34.965	ng	100
29) 2,4-Dichlorophenol	7.89	162	302978	38.608	ng	98
30) 1,2,4-Trichlorobenzene	7.97	180	336521	36.513	ng	100
31) Naphthalene	8.05	128	961621	37.765	ng	100
32) Benzoic acid	7.81	122	143400	32.722	ng	99
33) 4-Chloroaniline	8.10	127	153630	13.741	ng	98
34) Hexachlorobutadiene	8.17	225	203937	36.615	ng	99
35) Caprolactam	8.45	113	95004m	36.145	ng	
36) 4-Chloro-3-methylphenol	8.59	107	296271	38.547	ng	96
37) 2-Methylnaphthalene	8.74	142	669566	39.170	ng	99
38) 1-Methylnaphthalene	8.84	142	621592	38.485	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.91	216	351960	36.289	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.90	237	111062	33.296	nq	98
43) 2,4,6-Trichlorophenol	9.03	196	214815	35.980	nq	97
44) 2,4,5-Trichlorophenol	9.07	196	233656	37.905	nq	99
46) 1,1'-Biphenyl	9.20	154	828622	35.825	nq	99
47) 2-Chloronaphthalene	9.23	162	649187	36.298	nq	98
48) 2-Nitroaniline	9.33	65	150117	34.669	nq	95
49) Acenaphthylene	9.65	152	1028462	38.115	nq	99
50) Dimethylphthalate	9.51	163	1034744	48.598	nq	100
51) 2,6-Dinitrotoluene	9.57	165	174516	37.586	nq	91
52) Acenaphthene	9.82	154	591098	36.113	nq	100
53) 3-Nitroaniline	9.74	138	89309	16.787	nq	100
54) 2,4-Dinitrophenol	9.86	184	94178	60.786	nq	97
55) Dibenzofuran	10.00	168	940790	38.450	nq	97
56) 4-Nitrophenol	9.92	139	213518	62.862	nq	99
57) 2,4-Dinitrotoluene	9.99	165	230383	37.422	nq	94
58) Fluorene	10.34	166	722539	41.434	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.12	232	194502	37.326	nq	96
60) Diethylphthalate	10.22	149	753106	37.261	nq	99
61) 4-Chlorophenyl-phenylether	10.33	204	376554	40.156	nq	96
62) 4-Nitroaniline	10.36	138	162522	31.763	nq	98
63) Azobenzene	10.49	77	597496	38.504	nq	98
65) 4,6-Dinitro-2-methylphenol	10.40	198	85424	33.526	nq	91
66) n-Nitrosodiphenylamine	10.45	169	639985	38.230	nq	99
67) 4-Bromophenyl-phenylether	10.83	248	246293	38.066	nq	97
68) Hexachlorobenzene	10.90	284	265684	36.135	nq	# 91
69) Atrazine	10.99	200	203466	37.036	nq	99
70) Pentachlorophenol	11.10	266	196343	66.487	nq	99
71) Phenanthrene	11.31	178	1155605	40.032	nq	100
72) Anthracene	11.36	178	1129573	39.465	nq	99
73) Carbazole	11.52	167	1010454	38.084	nq	99
74) Di-n-butylphthalate	11.86	149	1226908	37.960	nq	99
75) Fluoranthene	12.52	202	1398097	43.123	nq	100
77) Benzidine	12.64	184	555787	39.665	nq	98
78) Pyrene	12.75	202	1395886	41.826	nq	100
80) Butylbenzylphthalate	13.37	149	533472	36.683	nq	100
81) Benzo(a)anthracene	13.96	228	1158601	39.243	nq	100
82) 3,3'-Dichlorobenzidine	13.92	252	350371	29.877	nq	99
83) Chrysene	13.99	228	1120962	38.684	nq	99
84) Bis(2-ethylhexyl)phthalate	13.95	149	720630	39.400	nq	99
85) Di-n-octyl phthalate	14.57	149	1285577	38.779	nq	99
86) Indeno(1,2,3-cd)pyrene	16.92	276	1369513	37.835	nq	99
88) Benzo(b)fluoranthene	15.02	252	1210028	38.601	nq	99
89) Benzo(k)fluoranthene	15.05	252	1144765	38.358	nq	100
90) Benzo(a)pyrene	15.38	252	1156494	39.577	nq	99
91) Dibenzo(a,h)anthracene	16.93	278	1115303	39.353	nq	100
92) Benzo(g,h,i)perylene	17.36	276	1152875	40.032	nq	99

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

