

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
 Data File : BP002598.D
 Acq On : 01 Jul 2020 11:04
 Operator : CG/JU
 Sample : L2819-08MSD
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BU-03-062620MSD

Manual Integrations
 APPROVED

mohammad
 7/1/2020 2:33:43 PM

Quant Time: Jul 01 11:38:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	97267	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	466393	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	243989	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	625695	20.00	ng	0.00
76) Chrysene-d12	21.10	240	904707	20.00	ng	0.00
86) Perylene-d12	23.29	264	1170604	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	851703	148.25	ng	0.00
7) Phenol-d6	6.77	99	1273182	158.78	ng	0.00
23) Nitrobenzene-d5	8.72	82	896351	99.03	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	465136	156.74	ng	0.00
45) 2-Fluorobiphenyl	12.84	172	1452616	87.63	ng	0.00
79) Terphenyl-d14	19.57	244	3613885	86.22	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.15	88	83983	34.109	ng	# 56
3) Pyridine	3.53	79	297762	40.803	ng	99
4) n-Nitrosodimethylamine	3.44	42	135267	43.667	ng	89
6) Aniline	6.90	93	457171	45.385	ng	98
8) 2-Chlorophenol	7.15	128	380894	59.687	ng	98
9) Benzaldehyde	6.72	77	123763	27.545	ng	98
10) Phenol	6.79	94	477753	59.084	ng	95
11) bis(2-Chloroethyl)ether	7.00	93	363619	54.789	ng	97
12) 1,3-Dichlorobenzene	7.46	146	345486	46.917	ng	99
13) 1,4-Dichlorobenzene	7.60	146	356972	47.583	ng	99
14) 1,2-Dichlorobenzene	7.92	146	350516	48.539	ng	99
15) Benzyl Alcohol	7.82	79	367223	60.949	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.11	45	503471	54.449	ng	99
17) 2-Methylphenol	8.03	107	350592	57.926	ng	98
18) Hexachloroethane	8.64	117	121089	44.721	ng	96
19) n-Nitroso-di-n-propylamine	8.37	70	312429	59.386	ng	99
20) 3+4-Methylphenols	8.36	107	482920	59.182	ng	98
22) Acetophenone	8.39	105	586914	49.718	ng	98
24) Nitrobenzene	8.76	77	471594	49.222	ng	99
25) Isophorone	9.28	82	951370	52.440	ng	98
26) 2-Nitrophenol	9.47	139	221361	50.659	ng	97
27) 2,4-Dimethylphenol	9.55	122	383967	57.991	ng	97
28) bis(2-Chloroethoxy)methane	9.77	93	552298	52.997	ng	99
29) 2,4-Dichlorophenol	10.00	162	425093	55.708	ng	96
30) 1,2,4-Trichlorobenzene	10.22	180	381198	44.373	ng	99
31) Naphthalene	10.39	128	1204923	48.803	ng	100
32) Benzoic acid	9.70	122	202805	40.630	ng	99
33) 4-Chloroaniline	10.51	127	286495	26.528	ng	99
34) Hexachlorobutadiene	10.70	225	199931	38.022	ng	97
35) Caprolactam	11.27	113	104451m	40.720	ng	
36) 4-Chloro-3-methylphenol	11.66	107	344713	41.216	ng	99
37) 2-Methylnaphthalene	12.02	142	630015	34.997	ng	99
38) 1-Methylnaphthalene	12.24	142	609507	35.950	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.40	216	319917	41.992	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.39	237	186365	48.676	nq	98
43) 2,4,6-Trichlorophenol	12.65	196	253740	49.223	nq	99
44) 2,4,5-Trichlorophenol	12.72	196	284826	47.998	nq	99
46) 1,1'-Biphenyl	13.05	154	835815	45.287	nq	99
47) 2-Chloronaphthalene	13.09	162	675295	44.949	nq	98
48) 2-Nitroaniline	13.29	65	257905	53.810	nq	99
49) Acenaphthylene	13.94	152	1135206	48.012	nq	98
50) Dimethylphthalate	13.69	163	1132346	59.089	nq	100
51) 2,6-Dinitrotoluene	13.80	165	218600	52.818	nq	98
52) Acenaphthene	14.29	154	674065	48.320	nq	100
53) 3-Nitroaniline	14.13	138	158466	35.164	nq	98
54) 2,4-Dinitrophenol	14.35	184	255665	97.661	nq	98
55) Dibenzofuran	14.63	168	1111624	49.172	nq	99
56) 4-Nitrophenol	14.47	139	365215	101.972	nq	99
57) 2,4-Dinitrotoluene	14.60	165	316292	56.810	nq	97
58) Fluorene	15.28	166	875056	50.824	nq	96
59) 2,3,4,6-Tetrachlorophenol	14.87	232	272457	56.836	nq	99
60) Diethylphthalate	15.07	149	945032	49.593	nq	99
61) 4-Chlorophenyl-phenylether	15.29	204	427842	46.337	nq	97
62) 4-Nitroaniline	15.30	138	241328	53.579	nq	99
63) Azobenzene	15.58	77	972463	52.778	nq	99
65) 4,6-Dinitro-2-methylphenol	15.37	198	190463	47.334	nq	98
66) n-Nitrosodiphenylamine	15.50	169	833771	45.647	nq	98
67) 4-Bromophenyl-phenylether	16.18	248	289781	41.213	nq	97
68) Hexachlorobenzene	16.30	284	376022	49.919	nq	95
69) Atrazine	16.47	200	296291	50.554	nq	99
70) Pentachlorophenol	16.65	266	445111	102.619	nq	99
71) Phenanthrene	17.03	178	1636813	48.621	nq	100
72) Anthracene	17.12	178	1675820	50.024	nq	99
73) Carbazole	17.39	167	1696158	52.389	nq	100
74) Di-n-butylphthalate	17.97	149	1897984	49.778	nq	100
75) Fluoranthene	19.02	202	2854201	70.043	nq	99
77) Benzidine	19.20	184	1491817	Below	Cal	100
78) Pyrene	19.37	202	2911483	47.279	nq	99
80) Butylbenzylphthalate	20.26	149	1290847	48.015	nq	98
81) Benzo(a)anthracene	21.08	228	2878582	48.668	nq	98
82) 3,3'-Dichlorobenzidine	21.02	252	866250	40.004	nq	97
83) Chrysene	21.13	228	2607869	45.981	nq	98
84) Bis(2-ethylhexyl)phthalate	21.03	149	1749983	45.331	nq	97
85) Di-n-octyl phthalate	21.90	149	3403463	49.915	nq	100
87) Indeno(1,2,3-cd)pyrene	25.53	276	3210528	38.091	nq	100
88) Benzo(b)fluoranthene	22.64	252	3592698	48.098	nq	98
89) Benzo(k)fluoranthene	22.69	252	3264945	46.854	nq	98
90) Benzo(a)pyrene	23.20	252	3325666	48.518	nq	99
91) Dibenzo(a,h)anthracene	25.53	278	2606319	37.843	nq	99
92) Benzo(g,h,i)perylene	26.19	276	2701035	39.213	nq	98

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

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