

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041019\
 Data File : BG040434.D
 Acq On : 11 Apr 2019 5:05
 Operator : JU/SJ
 Sample : PB118790BS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118790BS

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:46:48 PM

Quant Time: Apr 11 07:10:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	63719	20.00	ng	-0.03
21) Naphthalene-d8	10.86	136	264626	20.00	ng	-0.03
39) Acenaphthene-d10	14.68	164	169429	20.00	ng	-0.03
64) Phenanthrene-d10	17.42	188	451179	20.00	ng	-0.03
76) Chrysene-d12	21.70	240	442361	20.00	ng	-0.03
87) Perylene-d12	24.98	264	438249	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	530654	138.45	ng	-0.01
7) Phenol-d6	7.21	99	717106	135.38	ng	-0.01
23) Nitrobenzene-d5	9.22	82	432574	86.89	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	280778	153.34	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	960640	84.01	ng	-0.03
79) Terphenyl-d14	20.02	244	1624149	82.60	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	68151	37.813	ng	96
3) Pyridine	3.89	79	160631	33.102	ng	98
4) n-Nitrosodimethylamine	3.81	42	86670	38.612	ng	# 97
6) Aniline	7.37	93	204616	29.731	ng	97
8) 2-Chlorophenol	7.61	128	185419	41.325	ng	97
9) Benzaldehyde	7.19	77	86452	23.793	ng	98
10) Phenol	7.23	94	239621	41.676	ng	99
11) bis(2-Chloroethyl)ether	7.47	93	175263	38.058	ng	97
12) 1,3-Dichlorobenzene	7.93	146	185132	37.957	ng	97
13) 1,4-Dichlorobenzene	8.08	146	188211	38.744	ng	99
14) 1,2-Dichlorobenzene	8.40	146	178581	38.441	ng	98
15) Benzyl Alcohol	8.29	79	165426	38.777	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.57	45	329819	37.318	ng	99
17) 2-Methylphenol	8.48	107	168000	42.226	ng	98
18) Hexachloroethane	9.12	117	67818	36.702	ng	97
19) n-Nitroso-di-n-propylamine	8.85	70	137203	36.126	ng	96
20) 3+4-Methylphenols	8.81	107	225981	41.927	ng	98
22) Acetophenone	8.88	105	270329	40.952	ng	# 96
24) Nitrobenzene	9.27	77	206075	39.973	ng	98
25) Isophorone	9.78	82	396379	38.695	ng	98
26) 2-Nitrophenol	9.97	139	100990	41.341	ng	95
27) 2,4-Dimethylphenol	10.02	122	180618	46.548	ng	96
28) bis(2-Chloroethoxy)methane	10.26	93	233547	38.536	ng	99
29) 2,4-Dichlorophenol	10.51	162	185232	43.979	ng	99
30) 1,2,4-Trichlorobenzene	10.72	180	189353	40.944	ng	99
31) Naphthalene	10.91	128	555771	40.974	ng	99
32) Benzoic acid	10.17	122	79710	34.000	ng	92
33) 4-Chloroaniline	11.03	127	151231	26.138	ng	97
34) Hexachlorobutadiene	11.17	225	109939	37.170	ng	100
35) Caprolactam	11.82	113	71750m	42.928	ng	
36) 4-Chloro-3-methylphenol	12.14	107	216697	44.754	ng	96
37) 2-Methylnaphthalene	12.51	142	414500	41.319	ng	99
38) 1-Methylnaphthalene	12.73	142	395009	40.606	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	209503	38.905	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.84	237	229467	77.607	nq	98
43) 2,4,6-Trichlorophenol	13.11	196	155897	44.902	nq	96
44) 2,4,5-Trichlorophenol	13.19	196	168650	44.566	nq	98
46) 1,1'-Biphenyl	13.51	154	534275	39.910	nq	97
47) 2-Chloronaphthalene	13.56	162	415570	40.469	nq	99
48) 2-Nitroaniline	13.77	65	152901	44.143	nq	99
49) Acenaphthylene	14.40	152	702974	40.376	nq	99
50) Dimethylphthalate	14.13	163	559075	42.162	nq	100
51) 2,6-Dinitrotoluene	14.26	165	130952	43.982	nq	97
52) Acenaphthene	14.74	154	409134	41.010	nq	98
53) 3-Nitroaniline	14.59	138	104621	33.196	nq	94
54) 2,4-Dinitrophenol	14.81	184	128925	81.421	nq	97
55) Dibenzofuran	15.08	168	687662	42.896	nq	99
56) 4-Nitrophenol	14.90	139	221423	87.049	nq	99
57) 2,4-Dinitrotoluene	15.05	165	189930	45.909	nq	99
58) Fluorene	15.72	166	546869	43.390	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.30	232	167482	48.771	nq	99
60) Diethylphthalate	15.48	149	598362	44.098	nq	100
61) 4-Chlorophenyl-phenylether	15.71	204	289512	42.973	nq	96
62) 4-Nitroaniline	15.76	138	153766	45.463	nq	98
63) Azobenzene	16.00	77	558464	43.633	nq	98
65) 4,6-Dinitro-2-methylphenol	15.81	198	105482	41.614	nq	99
66) n-Nitrosodiphenylamine	15.93	169	527053	39.920	nq	96
67) 4-Bromophenyl-phenylether	16.60	248	197252	38.850	nq	98
68) Hexachlorobenzene	16.72	284	201577	39.486	nq	95
69) Atrazine	16.87	200	196444	39.998	nq	98
70) Pentachlorophenol	17.07	266	227901	77.217	nq	97
71) Phenanthrene	17.47	178	952984	40.185	nq	99
72) Anthracene	17.56	178	984025	41.456	nq	100
73) Carbazole	17.83	167	925510	41.200	nq	99
74) Di-n-butylphthalate	18.36	149	1111877	41.756	nq	99
75) Fluoranthene	19.47	202	1160856	41.339	nq	99
77) Benzidine	19.66	184	393020	24.604	nq	97
78) Pyrene	19.84	202	1149214	39.505	nq	97
80) Butylbenzylphthalate	20.71	149	514708	41.348	nq	96
81) Benzo(a)anthracene	21.67	228	1057751	38.821	nq	98
82) 3,3'-Dichlorobenzidine	21.59	252	297730	28.725	nq	98
83) Chrysene	21.75	228	1045906	39.941	nq	99
84) Bis(2-ethylhexyl)phthalate	21.55	149	747783	42.024	nq	100
85) Di-n-octyl phthalate	22.77	149	1268414	41.838	nq	100
86) Indeno(1,2,3-cd)pyrene	28.74	276	1290237	40.286	nq	# 91
88) Benzo(b)fluoranthene	23.93	252	1133410	42.242	nq	99
89) Benzo(k)fluoranthene	24.00	252	1101232	44.631	nq	99
90) Benzo(a)pyrene	24.82	252	1114735	44.280	nq	100
91) Dibenzo(a,h)anthracene	28.80	278	1057459	45.227	nq	99
92) Benzo(g,h,i)perylene	29.93	276	1078753	44.894	nq	98

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

