

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
 Data File : BM020109.D
 Acq On : 03 May 2019 13:28
 Operator : JU/SJ
 Sample : PB119431BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB119431BS

Manual Integrations
 APPROVED

mohammad
 5/7/2019 11:59:00 PM

Quant Time: May 04 04:03:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	182432	20.00	ng	0.00
21) Naphthalene-d8	10.58	136	681637	20.00	ng	0.00
39) Acenaphthene-d10	14.43	164	402394	20.00	ng	0.00
64) Phenanthrene-d10	17.17	188	897312	20.00	ng	0.00
76) Chrysene-d12	21.36	240	754674	20.00	ng	-0.01
87) Perylene-d12	23.64	264	604088	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	1557457	139.55	ng	0.00
7) Phenol-d6	6.96	99	2084271	136.70	ng	0.00
23) Nitrobenzene-d5	8.96	82	1298986	75.24	ng	0.00
42) 2,4,6-Tribromophenol	15.92	330	838449	134.24	ng	0.00
45) 2-Fluorobiphenyl	13.05	172	2546846	77.87	ng	0.00
79) Terphenyl-d14	19.80	244	3713772	85.83	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.31	88	167954	30.683	ng	95
3) Pyridine	3.70	79	426948	30.498	ng	99
4) n-Nitrosodimethylamine	3.62	42	170665	38.017	ng	80
6) Aniline	7.13	93	586609	30.563	ng	98
8) 2-Chlorophenol	7.35	128	473821	38.990	ng	98
9) Benzaldehyde	6.94	77	239201	20.728	ng	97
10) Phenol	6.98	94	639892	38.985	ng	97
11) bis(2-Chloroethyl)ether	7.22	93	447085	37.476	ng	98
12) 1,3-Dichlorobenzene	7.67	146	509695	36.049	ng	97
13) 1,4-Dichlorobenzene	7.82	146	524712	36.736	ng	95
14) 1,2-Dichlorobenzene	8.14	146	506153	37.195	ng	97
15) Benzyl Alcohol	8.03	79	544614	37.043	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.32	45	340294	34.400	ng	97
17) 2-Methylphenol	8.22	107	410652	38.865	ng	97
18) Hexachloroethane	8.86	117	213005	35.766	ng	97
19) n-Nitroso-di-n-propylamine	8.60	70	468930	35.948	ng	96
20) 3+4-Methylphenols	8.55	107	538843	38.371	ng	98
22) Acetophenone	8.62	105	759338	40.761	ng	# 97
24) Nitrobenzene	9.00	77	678529	37.904	ng	98
25) Isophorone	9.52	82	1140487	38.305	ng	99
26) 2-Nitrophenol	9.70	139	230873	42.719	ng	99
27) 2,4-Dimethylphenol	9.75	122	416305	46.269	ng	99
28) bis(2-Chloroethoxy)methane	10.00	93	629389	38.813	ng	99
29) 2,4-Dichlorophenol	10.23	162	469618	41.392	ng	97
30) 1,2,4-Trichlorobenzene	10.44	180	555927	39.822	ng	97
31) Naphthalene	10.63	128	1375929	38.898	ng	99
32) Benzoic acid	9.89	122	240871	39.055	ng	94
33) 4-Chloroaniline	10.75	127	328720	22.579	ng	96
34) Hexachlorobutadiene	10.90	225	379730	38.452	ng	95
35) Caprolactam	11.54	113	126352m	37.245	ng	
36) 4-Chloro-3-methylphenol	11.86	107	500826	37.564	ng	98
37) 2-Methylnaphthalene	12.25	142	1024465	39.727	ng	98
38) 1-Methylnaphthalene	12.46	142	964252	38.238	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.61	216	665046	42.243	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
 Data File : BM020109.D
 Acq On : 03 May 2019 13:28
 Operator : JU/SJ
 Sample : PB119431BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB119431BS

Manual Integrations
 APPROVED

mohammad
 5/7/2019 11:59:00 PM

Quant Time: May 04 04:03:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.58	237	804583	86.798	nq	99
43) 2,4,6-Trichlorophenol	12.85	196	395751	44.231	nq	97
44) 2,4,5-Trichlorophenol	12.92	196	425051	42.525	nq	96
46) 1,1'-Biphenyl	13.26	154	1330519	40.172	nq	99
47) 2-Chloronaphthalene	13.30	162	1068697	40.413	nq	99
48) 2-Nitroaniline	13.52	65	357916	38.017	nq	95
49) Acenaphthylene	14.15	152	1610498	38.054	nq	99
50) Dimethylphthalate	13.89	163	1279606	37.415	nq	99
51) 2,6-Dinitrotoluene	14.02	165	276355	38.365	nq	94
52) Acenaphthene	14.50	154	941884	38.809	nq	98
53) 3-Nitroaniline	14.34	138	173357	25.930	nq	97
54) 2,4-Dinitrophenol	14.55	184	323965	88.751	nq	95
55) Dibenzofuran	14.83	168	1565214	38.235	nq	98
56) 4-Nitrophenol	14.64	139	423968	74.326	nq	96
57) 2,4-Dinitrotoluene	14.80	165	369903	37.792	nq	94
58) Fluorene	15.48	166	1264338	37.606	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.05	232	389497	41.006	nq	97
60) Diethylphthalate	15.26	149	1254544	36.410	nq	99
61) 4-Chlorophenyl-phenylether	15.47	204	736736	38.675	nq	97
62) 4-Nitroaniline	15.51	138	243567	33.012	nq	97
63) Azobenzene	15.77	77	1459627	35.908	nq	99
65) 4,6-Dinitro-2-methylphenol	15.56	198	188612	42.605	nq	100
66) n-Nitrosodiphenylamine	15.69	169	1087080	41.171	nq	98
67) 4-Bromophenyl-phenylether	16.37	248	447857	40.702	nq	97
68) Hexachlorobenzene	16.47	284	504435	39.838	nq	97
69) Atrazine	16.64	200	378989	36.729	nq	99
70) Pentachlorophenol	16.82	266	619518	85.513	nq	97
71) Phenanthrene	17.22	178	1912732	38.616	nq	99
72) Anthracene	17.31	178	1930331	38.978	nq	99
73) Carbazole	17.58	167	1627472	36.420	nq	98
74) Di-n-butylphthalate	18.14	149	1946905	36.203	nq	98
75) Fluoranthene	19.23	202	2211932	35.294	nq	99
77) Benzidine	19.43	184	626632	25.972	nq	99
78) Pyrene	19.60	202	2218706	43.789	nq	100
80) Butylbenzylphthalate	20.50	149	781895	42.435	nq	99
81) Benzo(a)anthracene	21.34	228	1995424	37.703	nq	99
82) 3,3'-Dichlorobenzidine	21.28	252	698840	34.596	nq	100
83) Chrysene	21.40	228	1973554	38.450	nq	100
84) Bis(2-ethylhexyl)phthalate	21.27	149	1094421	38.278	nq	100
85) Di-n-octyl phthalate	22.16	149	1749483	36.815	nq	99
86) Indeno(1,2,3-cd)pyrene	25.97	276	2169015	35.526	nq	# 100
88) Benzo(b)fluoranthene	22.95	252	1958343	49.093	nq	100
89) Benzo(k)fluoranthene	23.00	252	1876156	51.560	nq	99
90) Benzo(a)pyrene	23.54	252	1830858	50.529	nq	99
91) Dibenzo(a,h)anthracene	25.98	278	1865577	49.509	nq	99
92) Benzo(g,h,i)perylene	26.69	276	1793706	50.139	nq	98

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

