

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092419\
 Data File : BM022751.D
 Acq On : 24 Sep 2019 15:48
 Operator : JU
 Sample : PB123301BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB123301BS

Manual Integrations
 APPROVED

mohammad
 9/26/2019 8:25:18 AM

Quant Time: Sep 25 04:09:00 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	250134	20.00	ng	0.00
21) Naphthalene-d8	10.62	136	1055769	20.00	ng	0.00
39) Acenaphthene-d10	14.46	164	658399	20.00	ng	0.00
64) Phenanthrene-d10	17.20	188	1355871	20.00	ng	0.00
76) Chrysene-d12	21.37	240	1350498	20.00	ng	0.00
87) Perylene-d12	23.64	264	1298143	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1525546	96.73	ng	0.00
7) Phenol-d6	6.99	99	2006948	99.07	ng	0.00
23) Nitrobenzene-d5	8.97	82	1069716	56.24	ng	0.00
42) 2,4,6-Tribromophenol	15.94	330	784611	95.46	ng	0.00
45) 2-Fluorobiphenyl	13.08	172	2630028	55.01	ng	0.00
79) Terphenyl-d14	19.82	244	4080183	58.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.31	88	217114	34.211	ng	# 100
3) Pyridine	3.71	79	532386	31.827	ng	98
4) n-Nitrosodimethylamine	3.62	42	227623	37.379	ng	# 96
6) Aniline	7.15	93	752658	30.987	ng	100
8) 2-Chlorophenol	7.39	128	656311	40.843	ng	99
9) Benzaldehyde	6.96	77	338803	29.373	ng	100
10) Phenol	7.02	94	792506	41.555	ng	99
11) bis(2-Chloroethyl)ether	7.25	93	583004	38.910	ng	100
12) 1,3-Dichlorobenzene	7.71	146	685358	35.685	ng	99
13) 1,4-Dichlorobenzene	7.86	146	701829	36.096	ng	100
14) 1,2-Dichlorobenzene	8.17	146	664188	35.831	ng	99
15) Benzyl Alcohol	8.06	79	520035	41.027	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.35	45	806292	37.044	ng	99
17) 2-Methylphenol	8.26	107	531119	42.075	ng	99
18) Hexachloroethane	8.90	117	249133	34.935	ng	96
19) n-Nitroso-di-n-propylamine	8.63	70	446693	39.489	ng	100
20) 3+4-Methylphenols	8.59	107	712932	42.403	ng	99
22) Acetophenone	8.64	105	904322	34.701	ng	# 99
24) Nitrobenzene	9.02	77	665647	36.579	ng	99
25) Isophorone	9.55	82	1222798	39.101	ng	99
26) 2-Nitrophenol	9.73	139	307667	39.164	ng	99
27) 2,4-Dimethylphenol	9.79	122	607762	46.212	ng	100
28) bis(2-Chloroethoxy)methane	10.03	93	807118	37.297	ng	100
29) 2,4-Dichlorophenol	10.26	162	609507	40.991	ng	99
30) 1,2,4-Trichlorobenzene	10.48	180	641470	35.064	ng	98
31) Naphthalene	10.66	128	1942899	37.245	ng	100
32) Benzoic acid	9.93	122	344304	38.920	ng	95
33) 4-Chloroaniline	10.77	127	510555	22.541	ng	99
34) Hexachlorobutadiene	10.96	225	364613	33.435	ng	99
35) Caprolactam	11.56	113	188426m	42.745	ng	
36) 4-Chloro-3-methylphenol	11.89	107	621216	42.822	ng	99
37) 2-Methylnaphthalene	12.27	142	1427343	39.308	ng	99
38) 1-Methylnaphthalene	12.50	142	1332005	38.645	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.65	216	735763	32.980	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092419\
 Data File : BM022751.D
 Acq On : 24 Sep 2019 15:48
 Operator : JU
 Sample : PB123301BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 PB123301BS

Manual Integrations
 APPROVED

mohammad
 9/26/2019 8:25:18 AM

Quant Time: Sep 25 04:09:00 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)
41) Hexachlorocyclopentadiene	12.64	237	959795	78.852	ng	99
43) 2,4,6-Trichlorophenol	12.89	196	484207	37.605	ng	99
44) 2,4,5-Trichlorophenol	12.96	196	524151	39.064	ng	99
46) 1,1'-Biphenyl	13.29	154	1835697	33.688	ng	99
47) 2-Chloronaphthalene	13.33	162	1395247	34.353	ng	99
48) 2-Nitroaniline	13.53	65	378052	35.747	ng	97
49) Acenaphthylene	14.18	152	2262078	37.611	ng	100
50) Dimethylphthalate	13.92	163	1744884	36.672	ng	100
51) 2,6-Dinitrotoluene	14.03	165	384272	38.610	ng	99
52) Acenaphthene	14.52	154	1361545	34.603	ng	100
53) 3-Nitroaniline	14.36	138	290416	25.347	ng	100
54) 2,4-Dinitrophenol	14.56	184	408987	76.208	ng	94
55) Dibenzofuran	14.86	168	2131265	37.102	ng	100
56) 4-Nitrophenol	14.66	139	720941	81.863	ng	99
57) 2,4-Dinitrotoluene	14.82	165	527333	40.444	ng	99
58) Fluorene	15.50	166	1737217	37.215	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.08	232	456758	39.032	ng	99
60) Diethylphthalate	15.28	149	1762821	38.103	ng	100
61) 4-Chlorophenyl-phenylether	15.50	204	873194	36.406	ng	98
62) 4-Nitroaniline	15.52	138	432076	35.969	ng	99
63) Azobenzene	15.79	77	1602158	38.348	ng	99
65) 4,6-Dinitro-2-methylphenol	15.57	198	275620	41.488	ng	99
66) n-Nitrosodiphenylamine	15.71	169	1532477	36.997	ng	100
67) 4-Bromophenyl-phenylether	16.39	248	535437	35.505	ng	98
68) Hexachlorobenzene	16.51	284	600461	34.143	ng	99
69) Atrazine	16.66	200	480094	35.654	ng	100
70) Pentachlorophenol	16.84	266	771329	82.239	ng	99
71) Phenanthrene	17.24	178	2763001	36.115	ng	100
72) Anthracene	17.33	178	2803753	37.814	ng	99
73) Carbazole	17.59	167	2564761	37.172	ng	99
74) Di-n-butylphthalate	18.16	149	3003271	38.388	ng	100
75) Fluoranthene	19.25	202	3227576	37.386	ng	99
77) Benzidine	19.43	184	1470640	39.695	ng	99
78) Pyrene	19.62	202	3361118	37.977	ng	99
80) Butylbenzylphthalate	20.52	149	1335267	39.697	ng	100
81) Benzo(a)anthracene	21.36	228	3204905	36.616	ng	99
82) 3,3'-Dichlorobenzidine	21.29	252	797537	25.694	ng	100
83) Chrysene	21.41	228	3062237	36.087	ng	99
84) Bis(2-ethylhexyl)phthalate	21.29	149	1988975	39.908	ng	100
85) Di-n-octyl phthalate	22.18	149	3249491	40.910	ng	99
86) Indeno(1,2,3-cd)pyrene	25.96	276	2962149	29.247	ng	# 100
88) Benzo(b)fluoranthene	22.96	252	3024880	38.485	ng	99
89) Benzo(k)fluoranthene	23.01	252	2981925	38.236	ng	100
90) Benzo(a)pyrene	23.55	252	2852009	39.379	ng	99
91) Dibenzo(a,h)anthracene	25.97	278	2501092	33.015	ng	99
92) Benzo(g,h,i)perylene	26.67	276	2328260	32.357	ng	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092419\
 Data File : BM022751.D
 Acq On : 24 Sep 2019 15:48
 Operator : JU
 Sample : PB123301BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB123301BS

Manual Integrations
 APPROVED

mohammad
 9/26/2019 8:25:18 AM

Quant Time: Sep 25 04:09:00 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

