

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM013020\
 Data File : BM024663.D
 Acq On : 30 Jan 2020 17:57
 Operator : CG/JU
 Sample : PB126427BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB126427BS

Manual Integrations
 APPROVED

mohammad
 2/3/2020 7:55:35 AM

Quant Time: Jan 30 23:48:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	174952	20.00	ng	-0.02
21) Naphthalene-d8	10.19	136	607300	20.00	ng	-0.02
39) Acenaphthene-d10	14.08	164	390161	20.00	ng	-0.01
64) Phenanthrene-d10	16.83	188	852863	20.00	ng	-0.01
76) Chrysene-d12	21.04	240	964829	20.00	ng	-0.01
87) Perylene-d12	23.16	264	1234796	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.08	112	1017676	110.62	ng	0.00
7) Phenol-d6	6.63	99	1246410	98.35	ng	-0.01
23) Nitrobenzene-d5	8.57	82	719713	58.12	ng	-0.02
42) 2,4,6-Tribromophenol	15.58	330	707774	111.52	ng	-0.01
45) 2-Fluorobiphenyl	12.69	172	1926764	67.14	ng	-0.02
79) Terphenyl-d14	19.49	244	3243174	62.75	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.06	88	128542	35.323	ng	91
3) Pyridine	3.44	79	312507	29.623	ng	95
4) n-Nitrosodimethylamine	3.35	42	137595	29.385	ng	80
6) Aniline	6.77	93	429577	27.845	ng	96
8) 2-Chlorophenol	7.00	128	410160	39.165	ng	90
9) Benzaldehyde	6.58	77	191541	25.579	ng	91
10) Phenol	6.66	94	446077	35.312	ng	95
11) bis(2-Chloroethyl)ether	6.87	93	339783	33.599	ng	95
12) 1,3-Dichlorobenzene	7.32	146	490338	38.113	ng	95
13) 1,4-Dichlorobenzene	7.46	146	498707	38.209	ng	98
14) 1,2-Dichlorobenzene	7.77	146	476123	37.792	ng	98
15) Benzyl Alcohol	7.67	79	325583	31.160	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.97	45	250621	27.071	ng	95
17) 2-Methylphenol	7.88	107	311511	34.993	ng	94
18) Hexachloroethane	8.49	117	174545	34.632	ng	93
19) n-Nitroso-di-n-propylamine	8.24	70	260030	27.328	ng	96
20) 3+4-Methylphenols	8.21	107	410706	33.329	ng	96
22) Acetophenone	8.24	105	556087	35.511	ng	98
24) Nitrobenzene	8.61	77	411984	34.717	ng	93
25) Isophorone	9.14	82	701110	33.321	ng	96
26) 2-Nitrophenol	9.32	139	227885	41.472	ng	89
27) 2,4-Dimethylphenol	9.40	122	344040	46.074	ng	94
28) bis(2-Chloroethoxy)methane	9.63	93	431161	33.544	ng	99
29) 2,4-Dichlorophenol	9.85	162	407018	40.345	ng	97
30) 1,2,4-Trichlorobenzene	10.06	180	499686	41.069	ng	99
31) Naphthalene	10.24	128	1209094	39.419	ng	99
32) Benzoic acid	9.53	122	234631m	34.350	ng	
33) 4-Chloroaniline	10.35	127	237851	17.691	ng	95
34) Hexachlorobutadiene	10.54	225	330716	39.320	ng	96
35) Caprolactam	11.12	113	111516m	34.455	ng	
36) 4-Chloro-3-methylphenol	11.50	107	359932	33.553	ng	91
37) 2-Methylnaphthalene	11.86	142	900477	38.041	ng	98
38) 1-Methylnaphthalene	12.09	142	845693	37.559	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.25	216	555520	39.894	ng	99

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41) Hexachlorocyclopentadiene	12.23	237	733234	93.943	nq	100
43) 2,4,6-Trichlorophenol	12.50	196	356054	40.181	nq	98
44) 2,4,5-Trichlorophenol	12.57	196	375000	41.438	nq	# 94
46) 1,1'-Biphenyl	12.90	154	1142295	38.687	nq	99
47) 2-Chloronaphthalene	12.94	162	935587	39.250	nq	97
48) 2-Nitroaniline	13.15	65	210139	28.905	nq	88
49) Acenaphthylene	13.80	152	1421493	39.836	nq	100
50) Dimethylphthalate	13.56	163	1115526	35.358	nq	100
51) 2,6-Dinitrotoluene	13.66	165	250487	37.344	nq	# 82
52) Acenaphthene	14.15	154	842757	38.060	nq	99
53) 3-Nitroaniline	13.99	138	153111	22.008	nq	92
54) 2,4-Dinitrophenol	14.20	184	297543	74.179	nq	# 87
55) Dibenzofuran	14.49	168	1400196	38.773	nq	97
56) 4-Nitrophenol	14.32	139	424421	78.045	nq	# 85
57) 2,4-Dinitrotoluene	14.46	165	345008	35.889	nq	# 86
58) Fluorene	15.14	166	1103729	36.369	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.72	232	352367	38.396	nq	96
60) Diethylphthalate	14.94	149	1054856	32.934	nq	98
61) 4-Chlorophenyl-phenylether	15.15	204	620624	35.472	nq	97
62) 4-Nitroaniline	15.16	138	230324	29.993	nq	88
63) Azobenzene	15.44	77	829314	30.354	nq	92
65) 4,6-Dinitro-2-methylphenol	15.23	198	213988	42.673	nq	85
66) n-Nitrosodiphenylamine	15.36	169	967694	39.500	nq	99
67) 4-Bromophenyl-phenylether	16.04	248	406644	38.341	nq	96
68) Hexachlorobenzene	16.15	284	479354	39.385	nq	95
69) Atrazine	16.33	200	381038	40.520	nq	97
70) Pentachlorophenol	16.49	266	600906	81.055	nq	99
71) Phenanthrene	16.87	178	1785896	38.669	nq	99
72) Anthracene	16.97	178	1817432	40.218	nq	99
73) Carbazole	17.24	167	1594150	38.751	nq	99
74) Di-n-butylphthalate	17.83	149	1749769	33.271	nq	99
75) Fluoranthene	18.90	202	2239992	38.860	nq	99
77) Benzidine	19.10	184	1387245	61.739	nq	99
78) Pyrene	19.27	202	2322686	36.517	nq	99
80) Butylbenzylphthalate	20.20	149	826770	30.977	nq	94
81) Benzo(a)anthracene	21.03	228	2541763	37.957	nq	100
82) 3,3'-Dichlorobenzidine	20.97	252	816817	33.908	nq	98
83) Chrysene	21.08	228	2428921	38.012	nq	100
84) Bis(2-ethylhexyl)phthalate	21.00	149	1218105	30.932	nq	# 98
85) Di-n-octyl phthalate	21.85	149	2191141	33.678	nq	# 96
86) Indeno(1,2,3-cd)pyrene	25.23	276	3563727	51.170	nq	98
88) Benzo(b)fluoranthene	22.54	252	2909332	36.328	nq	99
89) Benzo(k)fluoranthene	22.58	252	2814720	36.007	nq	100
90) Benzo(a)pyrene	23.07	252	2706209	37.167	nq	99
91) Dibenzo(a,h)anthracene	25.24	278	3030546	43.852	nq	98
92) Benzo(g,h,i)perylene	25.85	276	3037684	48.408	nq	98

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

