

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM013120\
 Data File : BM024687.D
 Acq On : 31 Jan 2020 22:10
 Operator : CG/JU
 Sample : L1004-18MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BU-01-012920MS

Manual Integrations
 APPROVED

mohammad
 2/3/2020 4:11:15 PM

Quant Time: Feb 01 02:29:23 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	202903	20.00	ng	-0.02
21) Naphthalene-d8	10.18	136	726007	20.00	ng	-0.02
39) Acenaphthene-d10	14.07	164	496887	20.00	ng	-0.02
64) Phenanthrene-d10	16.83	188	1161749	20.00	ng	-0.02
76) Chrysene-d12	21.04	240	1261719	20.00	ng	-0.02
87) Perylene-d12	23.14	264	1454407	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.07	112	1464310	137.25	ng	-0.01
7) Phenol-d6	6.63	99	1737922	118.25	ng	-0.02
23) Nitrobenzene-d5	8.56	82	1179521	79.68	ng	-0.02
42) 2,4,6-Tribromophenol	15.58	330	1238627	153.25	ng	-0.01
45) 2-Fluorobiphenyl	12.69	172	3363895	92.04	ng	-0.02
79) Terphenyl-d14	19.49	244	6648651	98.38	ng	-0.02

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.08	88	165164	39.134	ng	# 59
3) Pyridine	3.45	79	363997	29.751	ng	93
4) n-Nitrosodimethylamine	3.36	42	193760	35.679	ng	86
6) Aniline	6.77	93	214806	12.006	ng	95
8) 2-Chlorophenol	7.00	128	588021	48.414	ng	92
9) Benzaldehyde	6.57	77	248852	28.655	ng	93
10) Phenol	6.65	94	656806	44.831	ng	96
11) bis(2-Chloroethyl)ether	6.87	93	479708	40.900	ng	94
12) 1,3-Dichlorobenzene	7.32	146	646910	43.356	ng	96
13) 1,4-Dichlorobenzene	7.46	146	660890	43.660	ng	97
14) 1,2-Dichlorobenzene	7.77	146	629446	43.079	ng	98
15) Benzyl Alcohol	7.67	79	487380	40.220	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.96	45	353714	32.943	ng	95
17) 2-Methylphenol	7.88	107	454650	44.037	ng	96
18) Hexachloroethane	8.49	117	224987	38.491	ng	90
19) n-Nitroso-di-n-propylamine	8.23	70	383185	34.724	ng	95
20) 3+4-Methylphenols	8.20	107	615535	43.070	ng	97
22) Acetophenone	8.24	105	795257	42.481	ng	96
24) Nitrobenzene	8.60	77	584028	41.168	ng	93
25) Isophorone	9.13	82	1064357	42.314	ng	97
26) 2-Nitrophenol	9.31	139	335827	51.123	ng	90
27) 2,4-Dimethylphenol	9.39	122	510578	57.197	ng	95
28) bis(2-Chloroethoxy)methane	9.62	93	651254	42.382	ng	99
29) 2,4-Dichlorophenol	9.85	162	634700	52.627	ng	98
30) 1,2,4-Trichlorobenzene	10.05	180	701637	48.238	ng	96
31) Naphthalene	10.23	128	1695232	46.231	ng	99
32) Benzoic acid	9.55	122	371444m	45.488	ng	
33) 4-Chloroaniline	10.36	127	61701	3.839	ng	94
34) Hexachlorobutadiene	10.53	225	461542	45.902	ng	99
35) Caprolactam	11.13	113	155927m	40.299	ng	
36) 4-Chloro-3-methylphenol	11.49	107	591123	46.095	ng	91
37) 2-Methylnaphthalene	11.86	142	1335814	47.205	ng	99
38) 1-Methylnaphthalene	12.08	142	1267975	47.105	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.24	216	831170	46.869	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM013120\
 Data File : BM024687.D
 Acq On : 31 Jan 2020 22:10
 Operator : CG/JU
 Sample : L1004-18MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BU-01-012920MS

Manual Integrations
 APPROVED

mohammad
 2/3/2020 4:11:15 PM

Quant Time: Feb 01 02:29:23 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)
41) Hexachlorocyclopentadiene	12.23	237	926825	93.240	nq	99
43) 2,4,6-Trichlorophenol	12.49	196	579518	51.353	nq	97
44) 2,4,5-Trichlorophenol	12.56	196	616548	53.496	nq	97
46) 1,1'-Biphenyl	12.90	154	1770860	47.093	nq	98
47) 2-Chloronaphthalene	12.93	162	1425032	46.942	nq	98
48) 2-Nitroaniline	13.15	65	339431	36.660	nq	85
49) Acenaphthylene	13.79	152	2233133	49.139	nq	100
50) Dimethylphthalate	13.55	163	1976665	49.196	nq	100
51) 2,6-Dinitrotoluene	13.66	165	429972	50.334	nq #	83
52) Acenaphthene	14.14	154	1336671	47.400	nq	99
53) 3-Nitroaniline	13.99	138	97899	11.049	nq	91
54) 2,4-Dinitrophenol	14.20	184	511846	100.198	nq #	85
55) Dibenzofuran	14.48	168	2258659	49.111	nq	96
56) 4-Nitrophenol	14.32	139	731133	105.567	nq	86
57) 2,4-Dinitrotoluene	14.46	165	593426	48.471	nq #	86
58) Fluorene	15.13	166	1843189	47.690	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.72	232	607777	52.002	nq	94
60) Diethylphthalate	14.93	149	1874192	45.946	nq	99
61) 4-Chlorophenyl-phenylether	15.14	204	1035761	46.483	nq	98
62) 4-Nitroaniline	15.16	138	367534	37.581	nq	91
63) Azobenzene	15.43	77	1411166	40.557	nq	94
65) 4,6-Dinitro-2-methylphenol	15.22	198	373644	54.700	nq	88
66) n-Nitrosodiphenylamine	15.36	169	1657523	49.669	nq	100
67) 4-Bromophenyl-phenylether	16.03	248	705430	48.828	nq	95
68) Hexachlorobenzene	16.14	284	815302	49.177	nq	96
69) Atrazine	16.33	200	679502	53.047	nq	98
70) Pentachlorophenol	16.49	266	1097234	108.653	nq	98
71) Phenanthrene	16.87	178	3067004	48.752	nq	99
72) Anthracene	16.96	178	3128834	50.829	nq	100
73) Carbazole	17.23	167	2794088	49.861	nq	99
74) Di-n-butylphthalate	17.83	149	3299297	46.055	nq	99
75) Fluoranthene	18.90	202	3931635	50.072	nq	99
77) Benzidine	19.09	184	1153754	39.265	nq	99
78) Pyrene	19.26	202	4042615	48.602	nq	100
80) Butylbenzylphthalate	20.20	149	1544375	44.249	nq	89
81) Benzo(a)anthracene	21.03	228	4209277	48.067	nq	100
82) 3,3'-Dichlorobenzidine	20.97	252	692081	21.970	nq	99
83) Chrysene	21.07	228	4036192	48.303	nq	99
84) Bis(2-ethylhexyl)phthalate	21.00	149	2265823	43.998	nq	97
85) Di-n-octyl phthalate	21.84	149	3937063	46.274	nq #	96
86) Indeno(1,2,3-cd)pyrene	25.22	276	5193543	57.025	nq	99
88) Benzo(b)fluoranthene	22.53	252	4551180	48.248	nq	99
89) Benzo(k)fluoranthene	22.57	252	4411191	47.909	nq	99
90) Benzo(a)pyrene	23.06	252	4060943	47.351	nq	100
91) Dibenzo(a,h)anthracene	25.23	278	4405520	54.122	nq	99
92) Benzo(g,h,i)perylene	25.84	276	4331967	58.609	nq	98

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

