

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
 Data File : BF119574.D
 Acq On : 27 Mar 2020 18:28
 Operator : CG/JU
 Sample : PB127849BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127849BS

Manual Integrations
 APPROVED

mohammad
 3/31/2020 8:15:17 AM

Quant Time: Mar 28 01:29:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 27 07:52:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	252993	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1057216	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	573332	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	1018856	20.00	ng	0.00
76) Chrysene-d12	14.01	240	722651	20.00	ng	0.00
87) Perylene-d12	15.47	264	627742	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1978135	124.47	ng	0.01
7) Phenol-d6	6.46	99	2540541	125.14	ng	0.00
23) Nitrobenzene-d5	7.40	82	1646960	84.66	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	755993	127.81	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	3035114	78.43	ng	0.00
79) Terphenyl-d14	12.96	244	3403632	92.28	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.65	88	290238	36.977 ng	99
3) Pyridine	3.37	79	811751	37.539 ng	95
4) n-Nitrosodimethylamine	3.33	42	449833	42.658 ng	97
6) Aniline	6.49	93	971999	34.691 ng	99
8) 2-Chlorophenol	6.62	128	820172	49.137 ng	99
9) Benzaldehyde	6.38	77	495994	38.513 ng	97
10) Phenol	6.47	94	1044726	48.228 ng	96
11) bis(2-Chloroethyl)ether	6.57	93	776055	44.793 ng	97
12) 1,3-Dichlorobenzene	6.77	146	813499	45.031 ng	96
13) 1,4-Dichlorobenzene	6.85	146	811208	44.893 ng	97
14) 1,2-Dichlorobenzene	7.00	146	764377	45.256 ng	99
15) Benzyl Alcohol	6.97	79	674433	44.164 ng	99
16) 2,2'-oxybis(1-Chloropropan	7.11	45	1548241	44.304 ng	99
17) 2-Methylphenol	7.09	107	668191	43.692 ng	98
18) Hexachloroethane	7.35	117	308114	46.529 ng	96
19) n-Nitroso-di-n-propylamine	7.25	70	562130	45.938 ng	99
20) 3+4-Methylphenols	7.24	107	804983	42.949 ng	# 85
22) Acetophenone	7.25	105	1053547	41.705 ng	97
24) Nitrobenzene	7.42	77	894887	43.046 ng	100
25) Isophorone	7.66	82	1577669	39.981 ng	99
26) 2-Nitrophenol	7.73	139	432551	52.844 ng	91
27) 2,4-Dimethylphenol	7.77	122	733385	43.330 ng	97
28) bis(2-Chloroethoxy)methane	7.87	93	1019161	43.677 ng	99
29) 2,4-Dichlorophenol	7.97	162	695249	44.706 ng	98
30) 1,2,4-Trichlorobenzene	8.06	180	726826	42.260 ng	99
31) Naphthalene	8.14	128	2394033	44.003 ng	99
32) Benzoic acid	7.89	122	555925	51.062 ng	97
33) 4-Chloroaniline	8.19	127	522704	20.967 ng	97
34) Hexachlorobutadiene	8.26	225	435921	45.301 ng	99
35) Caprolactam	8.56	113	227900m	44.777 ng	
36) 4-Chloro-3-methylphenol	8.67	107	779160	45.840 ng	97
37) 2-Methylnaphthalene	8.83	142	1648576	46.171 ng	99
38) 1-Methylnaphthalene	8.93	142	1494468	44.220 ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	672629	40.026 ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
 Data File : BF119574.D
 Acq On : 27 Mar 2020 18:28
 Operator : CG/JU
 Sample : PB127849BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127849BS

Manual Integrations
 APPROVED

mohammad
 3/31/2020 8:15:17 AM

Quant Time: Mar 28 01:29:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 27 07:52:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	830024	86.879	nq	100
43) 2,4,6-Trichlorophenol	9.11	196	510326	42.436	nq	100
44) 2,4,5-Trichlorophenol	9.15	196	549013	40.753	nq	97
46) 1,1'-Biphenyl	9.30	154	1973336	42.260	nq	99
47) 2-Chloronaphthalene	9.32	162	1543711	42.205	nq	98
48) 2-Nitroaniline	9.42	65	582983	46.177	nq	100
49) Acenaphthylene	9.75	152	2429168	42.291	nq	99
50) Dimethylphthalate	9.60	163	1757104	43.147	nq	99
51) 2,6-Dinitrotoluene	9.66	165	410931	47.077	nq	95
52) Acenaphthene	9.92	154	1685189	47.062	nq	99
53) 3-Nitroaniline	9.83	138	299660	27.192	nq	96
54) 2,4-Dinitrophenol	9.94	184	417410	107.738	nq	96
55) Dibenzofuran	10.09	168	2093853	40.784	nq	98
56) 4-Nitrophenol	9.99	139	682658	78.525	nq	99
57) 2,4-Dinitrotoluene	10.07	165	534467	48.605	nq	96
58) Fluorene	10.43	166	1607098	42.331	nq	97
59) 2,3,4,6-Tetrachlorophenol	10.20	232	455678	45.251	nq	99
60) Diethylphthalate	10.31	149	1836514	44.433	nq	99
61) 4-Chlorophenyl-phenylether	10.42	204	783168	42.184	nq	99
62) 4-Nitroaniline	10.45	138	427092	40.415	nq	98
63) Azobenzene	10.59	77	1808340	43.446	nq	96
65) 4,6-Dinitro-2-methylphenol	10.47	198	270365	63.283	nq	88
66) n-Nitrosodiphenylamine	10.54	169	1561450	46.684	nq	99
67) 4-Bromophenyl-phenylether	10.92	248	542181	46.726	nq	92
68) Hexachlorobenzene	10.98	284	595314	50.128	nq	# 92
69) Atrazine	11.07	200	493711	53.842	nq	99
70) Pentachlorophenol	11.17	266	640968	92.047	nq	98
71) Phenanthrene	11.39	178	2302653	43.586	nq	100
72) Anthracene	11.44	178	2361150	43.974	nq	99
73) Carbazole	11.60	167	2198680	41.370	nq	100
74) Di-n-butylphthalate	11.93	149	2754203	48.445	nq	100
75) Fluoranthene	12.58	202	2650078	46.350	nq	99
77) Benzidine	12.70	184	1436748	46.979	nq	99
78) Pyrene	12.81	202	2589336	43.660	nq	100
80) Butylbenzylphthalate	13.43	149	1180223	49.202	nq	98
81) Benzo(a)anthracene	14.00	228	1947767	41.448	nq	99
82) 3,3'-Dichlorobenzidine	13.96	252	470998	25.420	nq	99
83) Chrysene	14.04	228	2042017	42.105	nq	100
84) Bis(2-ethylhexyl)phthalate	13.99	149	1547290	54.111	nq	99
85) Di-n-octyl phthalate	14.61	149	2840644	56.486	nq	100
86) Indeno(1,2,3-cd)pyrene	16.93	276	1884692	41.262	nq	99
88) Benzo(b)fluoranthene	15.04	252	2133995	50.891	nq	97
89) Benzo(k)fluoranthene	15.08	252	1821079	45.608	nq	99
90) Benzo(a)pyrene	15.41	252	1761173	46.215	nq	99
91) Dibenzo(a,h)anthracene	16.96	278	1555417	46.664	nq	100
92) Benzo(g,h,i)perylene	17.38	276	1577516	47.109	nq	96

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

