

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032420\
 Data File : BF119497.D
 Acq On : 24 Mar 2020 20:47
 Operator : CG/JU
 Sample : PB127731BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127731BS

Manual Integrations
 APPROVED

mohammad
 3/25/2020 12:29:42 PM

Quant Time: Mar 25 00:27:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 18 15:11:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	299828	20.00	ng	-0.01
21) Naphthalene-d8	8.13	136	1153509	20.00	ng	-0.02
39) Acenaphthene-d10	9.89	164	612613	20.00	ng	-0.01
64) Phenanthrene-d10	11.37	188	1194013	20.00	ng	-0.02
76) Chrysene-d12	14.02	240	859389	20.00	ng	-0.01
87) Perylene-d12	15.49	264	730328	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	2028174	107.68	ng	0.00
7) Phenol-d6	6.47	99	2605044	108.27	ng	0.00
23) Nitrobenzene-d5	7.41	82	1641595	77.34	ng	-0.01
42) 2,4,6-Tribromophenol	10.68	330	787955	124.67	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	3123860	75.54	ng	-0.01
79) Terphenyl-d14	12.97	244	3658574	83.41	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.68	88	282192	30.336	ng	98
3) Pyridine	3.42	79	788789	30.779	ng	99
4) n-Nitrosodimethylamine	3.36	42	432340	34.595	ng	95
6) Aniline	6.50	93	924128	27.830	ng	99
8) 2-Chlorophenol	6.63	128	838697	42.398	ng	99
9) Benzaldehyde	6.39	77	545786	35.759	ng	98
10) Phenol	6.48	94	1028135m	40.048	ng	
11) bis(2-Chloroethyl)ether	6.57	93	793647	38.653	ng	97
12) 1,3-Dichlorobenzene	6.79	146	837407	39.113	ng	98
13) 1,4-Dichlorobenzene	6.86	146	847951	39.596	ng	98
14) 1,2-Dichlorobenzene	7.02	146	793285	39.631	ng	99
15) Benzyl Alcohol	6.98	79	720773	39.826	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.12	45	1536300	37.095	ng	99
17) 2-Methylphenol	7.09	107	680494	37.546	ng	98
18) Hexachloroethane	7.36	117	323725	41.250	ng	98
19) n-Nitroso-di-n-propylamine	7.26	70	583266	40.220	ng	98
20) 3+4-Methylphenols	7.25	107	830421m	37.386	ng	
22) Acetophenone	7.25	105	1085795	39.394	ng	# 88
24) Nitrobenzene	7.43	77	872044	38.446	ng	96
25) Isophorone	7.67	82	1613884	37.485	ng	97
26) 2-Nitrophenol	7.75	139	436543	48.880	ng	98
27) 2,4-Dimethylphenol	7.78	122	752857	40.768	ng	99
28) bis(2-Chloroethoxy)methane	7.88	93	1030287	40.468	ng	99
29) 2,4-Dichlorophenol	7.99	162	701319	41.332	ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	738298	39.344	ng	99
31) Naphthalene	8.15	128	2347949	39.554	ng	99
32) Benzoic acid	7.90	122	583966	49.160	ng	99
33) 4-Chloroaniline	8.20	127	467478	17.187	ng	96
34) Hexachlorobutadiene	8.27	225	430037	40.959	ng	98
35) Caprolactam	8.57	113	225386m	40.586	ng	
36) 4-Chloro-3-methylphenol	8.68	107	760005	40.981	ng	99
37) 2-Methylnaphthalene	8.85	142	1604857	41.195	ng	99
38) 1-Methylnaphthalene	8.95	142	1529169	41.470	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.01	216	689724	38.412	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	849667	83.233	ng	99
43) 2,4,6-Trichlorophenol	9.12	196	534609	41.604	ng	99
44) 2,4,5-Trichlorophenol	9.16	196	540344	37.537	ng	96
46) 1,1'-Biphenyl	9.31	154	1948127	39.045	ng	98
47) 2-Chloronaphthalene	9.33	162	1514838	38.760	ng	99
48) 2-Nitroaniline	9.43	65	550230	40.788	ng	99
49) Acenaphthylene	9.75	152	2358116	38.422	ng	99
50) Dimethylphthalate	9.62	163	1762919	40.514	ng	100
51) 2,6-Dinitrotoluene	9.67	165	405120	43.435	ng	97
52) Acenaphthene	9.92	154	1630961	42.627	ng	99
53) 3-Nitroaniline	9.84	138	265649	22.560	ng	95
54) 2,4-Dinitrophenol	9.95	184	376598	92.104	ng	96
55) Dibenzofuran	10.10	168	2145685	39.114	ng	99
56) 4-Nitrophenol	10.00	139	714498	76.918	ng	86
57) 2,4-Dinitrotoluene	10.07	165	537833	45.775	ng	89
58) Fluorene	10.44	166	1649305	40.657	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.21	232	471808	43.849	ng	99
60) Diethylphthalate	10.32	149	1849874	41.886	ng	99
61) 4-Chlorophenyl-phenylether	10.43	204	809506	40.807	ng	97
62) 4-Nitroaniline	10.46	138	425922	37.720	ng	99
63) Azobenzene	10.59	77	1767702	39.747	ng	97
65) 4,6-Dinitro-2-methylphenol	10.49	198	256486	51.227	ng	89
66) n-Nitrosodiphenylamine	10.55	169	1527680	38.974	ng	98
67) 4-Bromophenyl-phenylether	10.92	248	535600	39.387	ng	100
68) Hexachlorobenzene	10.99	284	568946	40.880	ng	97
69) Atrazine	11.08	200	487392	45.356	ng	98
70) Pentachlorophenol	11.18	266	637931	78.172	ng	98
71) Phenanthrene	11.40	178	2417082	39.040	ng	100
72) Anthracene	11.46	178	2474189	39.320	ng	100
73) Carbazole	11.61	167	2321664	37.276	ng	99
74) Di-n-butylphthalate	11.95	149	2946189	44.220	ng	99
75) Fluoranthene	12.59	202	2674334	39.913	ng	100
77) Benzdine	12.71	184	1375839	37.829	ng	98
78) Pyrene	12.82	202	2688670	38.121	ng	99
80) Butylbenzylphthalate	13.44	149	1332637	46.717	ng	98
81) Benzo(a)anthracene	14.01	228	2074269	37.117	ng	99
82) 3,3'-Dichlorobenzidine	13.97	252	589479	26.752	ng	98
83) Chrysene	14.05	228	2174660	37.705	ng	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	1735809	51.045	ng	100
85) Di-n-octyl phthalate	14.62	149	3067579	51.293	ng	98
86) Indeno(1,2,3-cd)pyrene	16.96	276	1892055	34.832	ng	98
88) Benzo(b)fluoranthene	15.06	252	2029698	41.604	ng	98
89) Benzo(k)fluoranthene	15.09	252	1906525	41.041	ng	99
90) Benzo(a)pyrene	15.43	252	1724149	38.888	ng	98
91) Dibenzo(a,h)anthracene	16.97	278	1577943	40.691	ng	99
92) Benzo(g,h,i)perylene	17.40	276	1549948	39.784	ng	97

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

