

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090919\
 Data File : BF116898.D
 Acq On : 9 Sep 2019 15:58
 Operator : HP/JU
 Sample : PB122923BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122923BS

Manual Integrations
 APPROVED

mohammad
 9/12/2019 9:28:02 AM

Quant Time: Sep 09 16:47:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 09 16:37:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	83662	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	361681	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	190562	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	311725	20.00	ng	0.00
76) Chrysene-d12	13.99	240	226384	20.00	ng	0.00
87) Perylene-d12	15.44	264	165034	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	562800	108.98	ng	0.01
7) Phenol-d6	6.47	99	722582	106.56	ng	0.00
23) Nitrobenzene-d5	7.40	82	418516	66.06	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	159518	125.22	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	740469	65.55	ng	0.00
79) Terphenyl-d14	12.93	244	736539	60.19	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.69	88	85800	35.992	ng	98
3) Pyridine	3.43	79	189597	27.469	ng	98
4) n-Nitrosodimethylamine	3.36	42	84194	35.522	ng	91
6) Aniline	6.50	93	251307	26.757	ng	99
8) 2-Chlorophenol	6.63	128	249390	44.239	ng	99
9) Benzaldehyde	6.38	77	110146	24.655	ng	98
10) Phenol	6.49	94	311760	41.519	ng	99
11) bis(2-Chloroethyl)ether	6.57	93	232366	39.743	ng	96
12) 1,3-Dichlorobenzene	6.78	146	243659	38.243	ng	97
13) 1,4-Dichlorobenzene	6.85	146	251659	39.674	ng	99
14) 1,2-Dichlorobenzene	7.01	146	234687	39.923	ng	99
15) Benzyl Alcohol	6.98	79	226956	41.232	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.11	45	225810	32.918	ng	98
17) 2-Methylphenol	7.09	107	209858	42.525	ng	98
18) Hexachloroethane	7.35	117	95569	38.779	ng	97
19) n-Nitroso-di-n-propylamine	7.25	70	176459	39.521	ng	96
20) 3+4-Methylphenols	7.25	107	261939	45.650	ng	93
22) Acetophenone	7.25	105	359170	41.248	ng	# 97
24) Nitrobenzene	7.42	77	272024	42.216	ng	98
25) Isophorone	7.66	82	495162	38.570	ng	100
26) 2-Nitrophenol	7.73	139	125799	52.513	ng	97
27) 2,4-Dimethylphenol	7.77	122	215559	44.554	ng	98
28) bis(2-Chloroethoxy)methane	7.86	93	295014	37.604	ng	100
29) 2,4-Dichlorophenol	7.98	162	201786	43.399	ng	99
30) 1,2,4-Trichlorobenzene	8.06	180	199653	39.589	ng	96
31) Naphthalene	8.14	128	699958	40.699	ng	99
32) Benzoic acid	7.90	122	138205	50.591	ng	98
33) 4-Chloroaniline	8.19	127	200733	25.641	ng	99
34) Hexachlorobutadiene	8.26	225	110685	40.253	ng	96
35) Caprolactam	8.55	113	63596m	36.548	ng	
36) 4-Chloro-3-methylphenol	8.67	107	236975	44.345	ng	94
37) 2-Methylnaphthalene	8.83	142	480194	41.964	ng	98
38) 1-Methylnaphthalene	8.93	142	444348	41.135	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	186354	40.794	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	153998	58.378	ng	98
43) 2,4,6-Trichlorophenol	9.11	196	132283	42.244	ng	100
44) 2,4,5-Trichlorophenol	9.15	196	145827	44.856	ng	97
46) 1,1'-Biphenyl	9.29	154	562958	38.932	ng	98
47) 2-Chloronaphthalene	9.32	162	443752	38.746	ng	97
48) 2-Nitroaniline	9.41	65	141213	42.649	ng	89
49) Acenaphthylene	9.73	152	686756	39.668	ng	99
50) Dimethylphthalate	9.59	163	508099	38.620	ng	100
51) 2,6-Dinitrotoluene	9.65	165	111488	44.415	ng	93
52) Acenaphthene	9.90	154	402558	37.844	ng	98
53) 3-Nitroaniline	9.82	138	96383	30.640	ng	96
54) 2,4-Dinitrophenol	9.93	184	95666	104.312	ng	93
55) Dibenzofuran	10.07	168	605864	40.402	ng	98
56) 4-Nitrophenol	9.99	139	195530	90.652	ng	98
57) 2,4-Dinitrotoluene	10.06	165	148907	48.515	ng	95
58) Fluorene	10.42	166	484660	42.417	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.20	232	110529	47.108	ng	99
60) Diethylphthalate	10.29	149	506283	38.417	ng	98
61) 4-Chlorophenyl-phenylether	10.41	204	206646	41.300	ng	97
62) 4-Nitroaniline	10.44	138	130930	42.889	ng	97
63) Azobenzene	10.57	77	529859	38.559	ng	99
65) 4,6-Dinitro-2-methylphenol	10.47	198	64819	55.146	ng	99
66) n-Nitrosodiphenylamine	10.53	169	427061	39.603	ng	98
67) 4-Bromophenyl-phenylether	10.90	248	118670	37.474	ng	96
68) Hexachlorobenzene	10.97	284	127724	38.550	ng	# 88
69) Atrazine	11.05	200	112278	37.176	ng	98
70) Pentachlorophenol	11.16	266	144511	90.164	ng	93
71) Phenanthrene	11.38	178	686198	39.544	ng	99
72) Anthracene	11.43	178	700655	40.111	ng	100
73) Carbazole	11.59	167	669208	40.219	ng	100
74) Di-n-butylphthalate	11.91	149	809730	37.960	ng	99
75) Fluoranthene	12.56	202	711014	41.694	ng	99
77) Benzidine	12.68	184	148415	16.715	ng	98
78) Pyrene	12.79	202	721198	35.838	ng	99
80) Butylbenzylphthalate	13.40	149	345693	35.324	ng	99
81) Benzo(a)anthracene	13.98	228	581790	40.606	ng	100
82) 3,3'-Dichlorobenzidine	13.94	252	174219	35.981	ng	98
83) Chrysene	14.02	228	572280	38.769	ng	99
84) Bis(2-ethylhexyl)phthalate	13.96	149	471982	39.070	ng	99
85) Di-n-octyl phthalate	14.57	149	776374	39.272	ng	99
86) Indeno(1,2,3-cd)pyrene	16.89	276	407584	34.376	ng	99
88) Benzo(b)fluoranthene	15.02	252	525771	52.695	ng	98
89) Benzo(k)fluoranthene	15.05	252	457879	50.331	ng	97
90) Benzo(a)pyrene	15.38	252	450956	51.949	ng	99
91) Dibenzo(a,h)anthracene	16.90	278	344977	45.402	ng	98
92) Benzo(g,h,i)perylene	17.33	276	335697	43.320	ng	99

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

