

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030220\
 Data File : BF119170.D
 Acq On : 2 Mar 2020 15:47
 Operator : CG/JU
 Sample : PB127188BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127188BS

Manual Integrations
 APPROVED

mohammad
 3/3/2020 12:25:50 PM

Quant Time: Mar 03 07:48:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	146116	20.00	ng	-0.01
21) Naphthalene-d8	8.02	136	563240	20.00	ng	-0.02
39) Acenaphthene-d10	9.77	164	324619	20.00	ng	-0.02
64) Phenanthrene-d10	11.25	188	595634	20.00	ng	-0.01
76) Chrysene-d12	13.89	240	422647	20.00	ng	-0.01
87) Perylene-d12	15.32	264	275280	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	865344	98.97	ng	0.00
7) Phenol-d6	6.39	99	1066732	100.32	ng	-0.02
23) Nitrobenzene-d5	7.30	82	632941	59.33	ng	-0.02
42) 2,4,6-Tribromophenol	10.56	330	393566	92.68	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1315122	59.68	ng	-0.02
79) Terphenyl-d14	12.83	244	1551632	63.21	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.52	88	118488	30.438	ng	98
3) Pyridine	3.25	79	273167	25.694	ng	98
4) n-Nitrosodimethylamine	3.18	42	111258	34.546	ng	96
6) Aniline	6.40	93	301323	22.965	ng	# 83
8) 2-Chlorophenol	6.53	128	353924	37.509	ng	99
9) Benzaldehyde	6.29	77	87333	12.293	ng	99
10) Phenol	6.40	94	420476	36.574	ng	97
11) bis(2-Chloroethyl)ether	6.47	93	313667	35.658	ng	99
12) 1,3-Dichlorobenzene	6.67	146	383346	33.897	ng	98
13) 1,4-Dichlorobenzene	6.75	146	392867	34.715	ng	98
14) 1,2-Dichlorobenzene	6.90	146	366432	35.503	ng	99
15) Benzyl Alcohol	6.89	79	295555	38.458	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.01	45	256186	33.620	ng	# 57
17) 2-Methylphenol	7.00	107	289128	37.794	ng	99
18) Hexachloroethane	7.25	117	140085	34.599	ng	98
19) n-Nitroso-di-n-propylamine	7.15	70	237920	38.398	ng	98
20) 3+4-Methylphenols	7.16	107	367441	39.205	ng	96
22) Acetophenone	7.15	105	479850	36.821	ng	# 96
24) Nitrobenzene	7.32	77	370307	35.103	ng	99
25) Isophorone	7.56	82	663284	35.803	ng	100
26) 2-Nitrophenol	7.64	139	199765	35.740	ng	98
27) 2,4-Dimethylphenol	7.68	122	303800	40.049	ng	97
28) bis(2-Chloroethoxy)methane	7.77	93	406604	33.719	ng	99
29) 2,4-Dichlorophenol	7.89	162	315076	35.282	ng	99
30) 1,2,4-Trichlorobenzene	7.96	180	347996	32.867	ng	98
31) Naphthalene	8.04	128	1032567	35.349	ng	99
32) Benzoic acid	7.83	122	164529	32.268	ng	99
33) 4-Chloroaniline	8.09	127	257876	20.526	ng	97
34) Hexachlorobutadiene	8.16	225	209933	31.831	ng	98
35) Caprolactam	8.47	113	99449m	36.166	ng	
36) 4-Chloro-3-methylphenol	8.59	107	332469	36.786	ng	100
37) 2-Methylnaphthalene	8.73	142	716871	36.467	ng	98
38) 1-Methylnaphthalene	8.83	142	668270	36.148	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	356363	32.560	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	123171	36.681	nq	100
43) 2,4,6-Trichlorophenol	9.02	196	221694	32.076	nq	99
44) 2,4,5-Trichlorophenol	9.07	196	240255	33.126	nq	96
46) 1,1'-Biphenyl	9.19	154	868023	33.085	nq	98
47) 2-Chloronaphthalene	9.22	162	688603	32.570	nq	96
48) 2-Nitroaniline	9.32	65	197526	33.496	nq	97
49) Acenaphthylene	9.63	152	1047397	34.301	nq	99
50) Dimethylphthalate	9.50	163	845658	34.068	nq	99
51) 2,6-Dinitrotoluene	9.56	165	189460	34.354	nq	94
52) Acenaphthene	9.80	154	625585	33.976	nq	99
53) 3-Nitroaniline	9.73	138	130153	21.288	nq	100
54) 2,4-Dinitrophenol	9.85	184	149416	58.011	nq	# 89
55) Dibenzofuran	9.98	168	948629	34.518	nq	96
56) 4-Nitrophenol	9.92	139	196451	53.479	nq	94
57) 2,4-Dinitrotoluene	9.97	165	242438	33.915	nq	98
58) Fluorene	10.32	166	754273	35.320	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.10	232	201221	32.880	nq	98
60) Diethylphthalate	10.20	149	819090	35.015	nq	100
61) 4-Chlorophenyl-phenylether	10.31	204	396780	35.102	nq	96
62) 4-Nitroaniline	10.35	138	178830	28.589	nq	99
63) Azobenzene	10.47	77	722315	35.524	nq	95
65) 4,6-Dinitro-2-methylphenol	10.39	198	128092	35.539	nq	96
66) n-Nitrosodiphenylamine	10.43	169	681879	34.962	nq	99
67) 4-Bromophenyl-phenylether	10.80	248	258895	33.253	nq	95
68) Hexachlorobenzene	10.87	284	291243	32.445	nq	97
69) Atrazine	10.96	200	234049	34.347	nq	98
70) Pentachlorophenol	11.07	266	219936	60.824	nq	98
71) Phenanthrene	11.28	178	1112974	34.263	nq	98
72) Anthracene	11.33	178	1135854	34.997	nq	97
73) Carbazole	11.49	167	985737	34.092	nq	99
74) Di-n-butylphthalate	11.82	149	1248229	35.648	nq	98
75) Fluoranthene	12.46	202	1175383	34.089	nq	98
77) Benzidine	12.59	184	193350	11.249	nq	98
78) Pyrene	12.69	202	1175624	34.640	nq	99
80) Butylbenzylphthalate	13.30	149	508224	35.581	nq	96
81) Benzo(a)anthracene	13.88	228	980333	34.042	nq	99
82) 3,3'-Dichlorobenzidine	13.84	252	300510	26.874	nq	98
83) Chrysene	13.92	228	966416	33.679	nq	98
84) Bis(2-ethylhexyl)phthalate	13.86	149	706180	39.134	nq	98
85) Di-n-octyl phthalate	14.47	149	1102455	38.344	nq	99
86) Indeno(1,2,3-cd)pyrene	16.72	276	744694	26.803	nq	99
88) Benzo(b)fluoranthene	14.91	252	843603	46.492	nq	98
89) Benzo(k)fluoranthene	14.94	252	809942	48.167	nq	99
90) Benzo(a)pyrene	15.26	252	693798	42.366	nq	98
91) Dibenzo(a,h)anthracene	16.72	278	640312	44.438	nq	99
92) Benzo(g,h,i)perylene	17.13	276	612886	43.049	nq	99

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

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