

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121322\
 Data File : BF131612.D
 Acq On : 13 Dec 2022 14:49
 Operator : CG\JU
 Sample : PB149584BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB149584BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/14/2022
 Supervised By :mohammad ahmed 12/14/2022

Quant Time: Dec 14 03:32:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 14 03:22:29 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	71010	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	255066	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	149637	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	288944	20.000	ng	0.00
76) Chrysene-d12	14.092	240	175571	20.000	ng	0.00
86) Perylene-d12	15.592	264	146631	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	395461	92.908	ng	0.01
7) Phenol-d6	6.522	99	513684	92.094	ng	0.00
23) Nitrobenzene-d5	7.457	82	533325	79.051	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	145253	87.370	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	910519	76.722	ng	0.00
79) Terphenyl-d14	13.027	244	1074242	83.910	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.728	88	71363	37.144	ng	98
3) Pyridine	3.475	79	196653	36.867	ng	97
4) n-Nitrosodimethylamine	3.452	42	114328	38.863	ng	# 84
6) Aniline	6.545	93	254294	38.120	ng	92
8) 2-Chlorophenol	6.669	128	195273	43.291	ng	97
9) Benzaldehyde	6.434	77	127695	33.419	ng	96
10) Phenol	6.540	94	282679m	45.435	ng	
11) bis(2-Chloroethyl)ether	6.622	93	200287	42.945	ng	97
12) 1,3-Dichlorobenzene	6.822	146	213954	40.911	ng	98
13) 1,4-Dichlorobenzene	6.904	146	220128	41.730	ng	98
14) 1,2-Dichlorobenzene	7.051	146	206524	40.984	ng	98
15) Benzyl Alcohol	7.034	79	211022	41.267	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.157	45	242356	40.229	ng	95
17) 2-Methylphenol	7.145	107	174439	42.910	ng	97
18) Hexachloroethane	7.398	117	86209	40.513	ng	94
19) n-Nitroso-di-n-propyla...	7.304	70	164860	39.417	ng	96
20) 3+4-Methylphenols	7.298	107	225738	40.768	ng	# 80
22) Acetophenone	7.298	105	320045	40.040	ng	95
24) Nitrobenzene	7.475	77	260946	38.349	ng	94
25) Isophorone	7.716	82	448079	41.233	ng	99
26) 2-Nitrophenol	7.787	139	102728	44.271	ng	99
27) 2,4-Dimethylphenol	7.828	122	171324	48.170	ng	99
28) bis(2-Chloroethoxy)met...	7.922	93	245117	43.603	ng	98
29) 2,4-Dichlorophenol	8.034	162	175167	39.832	ng	96
30) 1,2,4-Trichlorobenzene	8.116	180	203875	37.803	ng	97
31) Naphthalene	8.198	128	559616	38.954	ng	100
32) Benzoic acid	7.963	122	113905	44.221	ng	97
33) 4-Chloroaniline	8.245	127	160560	29.816	ng	99
34) Hexachlorobutadiene	8.310	225	134851	35.479	ng	99
35) Caprolactam	8.628	113	54037m	47.501	ng	
36) 4-Chloro-3-methylphenol	8.734	107	208592	42.240	ng	98
37) 2-Methylnaphthalene	8.892	142	374255	38.093	ng	98
38) 1-Methylnaphthalene	8.992	142	355255	37.667	ng	98
40) 1,2,4,5-Tetrachloroben...	9.057	216	215822	38.962	ng	98
41) Hexachlorocyclopentadiene	9.039	237	244790	87.065	ng	99
43) 2,4,6-Trichlorophenol	9.175	196	139687	40.770	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.216	196	145433	39.470	ng	97
46) 1,1'-Biphenyl	9.357	154	475698	40.207	ng	100
47) 2-Chloronaphthalene	9.386	162	385883	40.032	ng	99
48) 2-Nitroaniline	9.486	65	140206	41.060	ng	91
49) Acenaphthylene	9.804	152	578079	40.844	ng	99
50) Dimethylphthalate	9.663	163	448234	39.294	ng	99
51) 2,6-Dinitrotoluene	9.728	165	99874	42.410	ng	85
52) Acenaphthene	9.975	154	426758m	45.009	ng	
53) 3-Nitroaniline	9.898	138	72476	32.076	ng	100
54) 2,4-Dinitrophenol	10.010	184	120349	88.799	ng	# 88
55) Dibenzofuran	10.151	168	533574	39.828	ng	97
56) 4-Nitrophenol	10.075	139	152642	97.247	ng	# 77
57) 2,4-Dinitrotoluene	10.133	165	138045	44.190	ng	98
58) Fluorene	10.492	166	426252	39.138	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.269	232	121151m	38.828	ng	
60) Diethylphthalate	10.369	149	414770	37.890	ng	99
61) 4-Chlorophenyl-phenyle...	10.481	204	227725	37.836	ng	98
62) 4-Nitroaniline	10.522	138	97503	43.547	ng	91
63) Azobenzene	10.645	77	466559	37.369	ng	97
65) 4,6-Dinitro-2-methylph...	10.545	198	76547	41.827	ng	83
66) n-Nitrosodiphenylamine	10.604	169	358536	38.382	ng	98
67) 4-Bromophenyl-phenylether	10.975	248	135013	36.323	ng	96
68) Hexachlorobenzene	11.039	284	153518	38.183	ng	95
69) Atrazine	11.133	200	134662	45.396	ng	97
70) Pentachlorophenol	11.239	266	159829	77.218	ng	98
71) Phenanthrene	11.463	178	636515	39.522	ng	100
72) Anthracene	11.516	178	633771	40.447	ng	100
73) Carbazole	11.669	167	555790	43.456	ng	99
74) Di-n-butylphthalate	11.998	149	612372	39.028	ng	99
75) Fluoranthene	12.657	202	696367	41.990	ng	99
77) Benzidine	12.780	184	192355	40.890	ng	99
78) Pyrene	12.886	202	707426	40.863	ng	99
80) Butylbenzylphthalate	13.504	149	220679	44.047	ng	97
81) Benzo(a)anthracene	14.080	228	512968	40.458	ng	99
82) 3,3'-Dichlorobenzidine	14.039	252	133650	37.049	ng	95
83) Chrysene	14.116	228	472143	39.169	ng	99
84) Bis(2-ethylhexyl)phtha...	14.063	149	249166	37.232	ng	98
85) Di-n-octyl phthalate	14.680	149	327590	31.325	ng	98
87) Indeno(1,2,3-cd)pyrene	17.127	276	447640	37.407	ng	99
88) Benzo(b)fluoranthene	15.151	252	424287	43.903	ng	99
89) Benzo(k)fluoranthene	15.180	252	416193	41.260	ng	99
90) Benzo(a)pyrene	15.533	252	370494	44.922	ng	100
91) Dibenzo(a,h)anthracene	17.145	278	382551	38.843	ng	99
92) Benzo(g,h,i)perylene	17.598	276	376238	38.010	ng	98

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

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