

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092923\
 Data File : BF135514.D
 Acq On : 29 Sep 2023 14:09
 Operator : CG\JU
 Sample : 04620-10MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

STONE-INMS

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 10/02/2023

Supervised By :mohammad ahmed 10/03/2023

Quant Time: Sep 30 00:19:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 30 00:12:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.746	152	153108	20.000	ng	0.00
21) Naphthalene-d8	8.028	136	562162	20.000	ng	# 0.00
39) Acenaphthene-d10	9.781	164	257862	20.000	ng	0.00
64) Phenanthrene-d10	11.275	188	411269	20.000	ng	# 0.00
76) Chrysene-d12	13.933	240	254064	20.000	ng	0.00
86) Perylene-d12	15.392	264	304026	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	1143260	121.447	ng	0.02
7) Phenol-d6	6.404	99	1320120	110.286	ng	0.00
23) Nitrobenzene-d5	7.316	82	949894	91.801	ng	0.00
42) 2,4,6-Tribromophenol	10.581	330	262673	102.506	ng	0.00
45) 2-Fluorobiphenyl	9.104	172	1504629	88.034	ng	0.00
79) Terphenyl-d14	12.869	244	1204435	73.490	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.634	88	200070	48.481	ng	# 78
3) Pyridine	3.375	79	392695	35.357	ng	96
4) n-Nitrosodimethylamine	3.305	42	294026	49.887	ng	94
6) Aniline	6.422	93	309070	19.690	ng	# 50
8) 2-Chlorophenol	6.540	128	478599	47.626	ng	98
9) Benzaldehyde	6.310	77	82058	12.034	ng	99
10) Phenol	6.416	94	511190	38.946	ng	97
11) bis(2-Chloroethyl)ether	6.493	93	488578	49.624	ng	98
12) 1,3-Dichlorobenzene	6.687	146	464659	45.498	ng	98
13) 1,4-Dichlorobenzene	6.769	146	469819	45.216	ng	99
14) 1,2-Dichlorobenzene	6.916	146	436398	45.226	ng	98
15) Benzyl Alcohol	6.898	79	405049	47.156	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.028	45	864498	46.583	ng	84
17) 2-Methylphenol	7.016	107	362350	40.861	ng	94
18) Hexachloroethane	7.251	117	164508	42.849	ng	92
19) n-Nitroso-di-n-propyla...	7.169	70	314284	42.389	ng	98
20) 3+4-Methylphenols	7.169	107	441594	41.412	ng	# 83
22) Acetophenone	7.163	105	620739	48.297	ng	93
24) Nitrobenzene	7.334	77	505288	49.406	ng	98
25) Isophorone	7.575	82	928147	48.849	ng	99
26) 2-Nitrophenol	7.651	139	229935	46.844	ng	91
27) 2,4-Dimethylphenol	7.693	122	367884	44.589	ng	96
28) bis(2-Chloroethoxy)met...	7.781	93	580789	51.071	ng	99
29) 2,4-Dichlorophenol	7.893	162	360706	47.181	ng	100
30) 1,2,4-Trichlorobenzene	7.969	180	385073	48.189	ng	99
31) Naphthalene	8.045	128	1288686	47.181	ng	99
32) Benzoic acid	7.828	122	237718	38.263	ng	99
33) 4-Chloroaniline	8.104	127	165177	13.783	ng	96
34) Hexachlorobutadiene	8.157	225	214552	44.528	ng	99
35) Caprolactam	8.481	113	95171m	37.092	ng	
36) 4-Chloro-3-methylphenol	8.598	107	374456	44.069	ng	99
37) 2-Methylnaphthalene	8.739	142	788639	42.954	ng	99
38) 1-Methylnaphthalene	8.839	142	760189	44.224	ng	98
40) 1,2,4,5-Tetrachloroben...	8.904	216	366488	49.058	ng	99
41) Hexachlorocyclopentadiene	8.887	237	131111	43.538	ng	98
43) 2,4,6-Trichlorophenol	9.028	196	236959	48.446	ng	99

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QLast Update : Sat Sep 30 00:12:20 2023

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.075	196	245739	43.626	ng	99
46) 1,1'-Biphenyl	9.204	154	974325	49.168	ng	98
47) 2-Chloronaphthalene	9.234	162	726731	50.545	ng	99
48) 2-Nitroaniline	9.339	65	277138	49.615	ng	96
49) Acenaphthylene	9.645	152	1152198	48.219	ng	99
50) Dimethylphthalate	9.510	163	827256	47.450	ng	100
51) 2,6-Dinitrotoluene	9.581	165	174159	45.600	ng	90
52) Acenaphthene	9.816	154	671497	47.570	ng	95
53) 3-Nitroaniline	9.751	138	130153	29.032	ng	92
54) 2,4-Dinitrophenol	9.869	184	102726	49.927	ng	95
55) Dibenzofuran	9.992	168	958657	44.505	ng	99
56) 4-Nitrophenol	9.934	139	180461	63.336	ng	95
57) 2,4-Dinitrotoluene	9.986	165	189112	39.552	ng	# 80
58) Fluorene	10.334	166	736232	44.460	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.116	232	180954	41.762	ng	99
60) Diethylphthalate	10.210	149	796913	45.547	ng	99
61) 4-Chlorophenyl-phenyle...	10.328	204	343051	43.126	ng	93
62) 4-Nitroaniline	10.369	138	173177	40.186	ng	95
63) Azobenzene	10.486	77	816607	47.688	ng	97
65) 4,6-Dinitro-2-methylph...	10.404	198	76247	34.906	ng	89
66) n-Nitrosodiphenylamine	10.451	169	663601	53.297	ng	100
67) 4-Bromophenyl-phenylether	10.816	248	209920	51.321	ng	93
68) Hexachlorobenzene	10.881	284	215815	51.790	ng	# 87
69) Atrazine	10.981	200	184182	54.976	ng	98
70) Pentachlorophenol	11.086	266	180696	72.475	ng	99
71) Phenanthrene	11.298	178	979380	50.350	ng	99
72) Anthracene	11.351	178	996110	49.821	ng	99
73) Carbazole	11.516	167	912423	50.131	ng	98
74) Di-n-butylphthalate	11.845	149	1156246	53.912	ng	100
75) Fluoranthene	12.492	202	991649	48.927	ng	98
77) Benzidine	12.622	184	159620	30.667	ng	99
78) Pyrene	12.722	202	1000211	41.350	ng	99
80) Butylbenzylphthalate	13.345	149	427538	50.201	ng	94
81) Benzo(a)anthracene	13.922	228	818167	49.871	ng	98
82) 3,3'-Dichlorobenzidine	13.886	252	167594	32.705	ng	# 97
83) Chrysene	13.957	228	810148	49.669	ng	100
84) Bis(2-ethylhexyl)phtha...	13.910	149	575187	65.815	ng	99
85) Di-n-octyl phthalate	14.527	149	1044755	75.224	ng	100
87) Indeno(1,2,3-cd)pyrene	16.874	276	846262	42.453	ng	96
88) Benzo(b)fluoranthene	14.969	252	887483	53.924	ng	# 98
89) Benzo(k)fluoranthene	14.998	252	730180	44.913	ng	# 97
90) Benzo(a)pyrene	15.333	252	735934	46.890	ng	99
91) Dibenzo(a,h)anthracene	16.886	278	696281	42.690	ng	98
92) Benzo(g,h,i)perylene	17.315	276	668672	39.255	ng	96

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

