

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110123\
 Data File : BF136070.D
 Acq On : 01 Nov 2023 11:51
 Operator : CG\JU
 Sample : PB156691BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB156691BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/02/2023
 Supervised By :mohammad ahmed 11/02/2023

Quant Time: Nov 01 22:37:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 31 01:13:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	143089	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	591757	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	309384	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	565792	20.000	ng	# 0.00
76) Chrysene-d12	14.021	240	290118	20.000	ng	0.00
86) Perylene-d12	15.510	264	274421	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	1187640	130.427	ng	0.02
7) Phenol-d6	6.469	99	1492827	131.581	ng	0.01
23) Nitrobenzene-d5	7.393	82	904725	92.165	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	460010	139.300	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1655354	85.292	ng	0.00
79) Terphenyl-d14	12.963	244	1868875	100.107	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.669	88	163300	41.184	ng	99
3) Pyridine	3.422	79	365010	33.726	ng	99
4) n-Nitrosodimethylamine	3.393	42	201994	44.639	ng	97
6) Aniline	6.493	93	522947	35.602	ng	98
8) 2-Chlorophenol	6.610	128	478670	49.169	ng	99
9) Benzaldehyde	6.375	77	137602	21.751	ng	96
10) Phenol	6.481	94	556734m	47.444	ng	
11) bis(2-Chloroethyl)ether	6.569	93	438814	47.767	ng	99
12) 1,3-Dichlorobenzene	6.769	146	484964	47.863	ng	98
13) 1,4-Dichlorobenzene	6.845	146	490986	48.353	ng	98
14) 1,2-Dichlorobenzene	6.998	146	470115	49.513	ng	99
15) Benzyl Alcohol	6.969	79	388931	48.348	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.104	45	611848	48.388	ng	99
17) 2-Methylphenol	7.081	107	373825	45.265	ng	97
18) Hexachloroethane	7.340	117	175236	49.100	ng	92
19) n-Nitroso-di-n-propyla...	7.245	70	302961	46.736	ng	97
20) 3+4-Methylphenols	7.234	107	444836	44.172	ng	93
22) Acetophenone	7.240	105	607743	46.315	ng	# 91
24) Nitrobenzene	7.410	77	456085	45.932	ng	95
25) Isophorone	7.651	82	865885	46.723	ng	99
26) 2-Nitrophenol	7.728	139	239988	52.974	ng	96
27) 2,4-Dimethylphenol	7.763	122	410371	47.936	ng	99
28) bis(2-Chloroethoxy)met...	7.863	93	532131	46.864	ng	99
29) 2,4-Dichlorophenol	7.969	162	387437	48.007	ng	98
30) 1,2,4-Trichlorobenzene	8.051	180	417306	47.400	ng	99
31) Naphthalene	8.134	128	1315098	46.086	ng	99
32) Benzoic acid	7.892	122	324638	47.408	ng	97
33) 4-Chloroaniline	8.181	127	394543	31.572	ng	97
34) Hexachlorobutadiene	8.245	225	241709	47.873	ng	99
35) Caprolactam	8.557	113	134240m	51.154	ng	
36) 4-Chloro-3-methylphenol	8.663	107	406134	47.934	ng	99
37) 2-Methylnaphthalene	8.822	142	862477	45.076	ng	99
38) 1-Methylnaphthalene	8.922	142	802672	45.079	ng	99
40) 1,2,4,5-Tetrachloroben...	8.992	216	405066	45.908	ng	99
41) Hexachlorocyclopentadiene	8.975	237	480913	102.021	ng	99
43) 2,4,6-Trichlorophenol	9.098	196	278462	48.403	ng	97

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44) 2,4,5-Trichlorophenol	9.145	196	303008	45.469	ng	98
46) 1,1'-Biphenyl	9.292	154	1079382	45.766	ng	99
47) 2-Chloronaphthalene	9.316	162	813896	46.811	ng	99
48) 2-Nitroaniline	9.410	65	249429	48.475	ng	97
49) Acenaphthylene	9.734	152	1248463	46.028	ng	99
50) Dimethylphthalate	9.598	163	952993	46.698	ng	100
51) 2,6-Dinitrotoluene	9.657	165	217748	49.888	ng	97
52) Acenaphthene	9.904	154	786912	45.637	ng	100
53) 3-Nitroaniline	9.822	138	203012	39.862	ng	93
54) 2,4-Dinitrophenol	9.934	184	225781	93.905	ng	96
55) Dibenzofuran	10.075	168	1144169	45.507	ng	97
56) 4-Nitrophenol	9.986	139	365917	96.053	ng	87
57) 2,4-Dinitrotoluene	10.063	165	282931	51.233	ng	94
58) Fluorene	10.422	166	853000	45.312	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.192	232	256267	48.369	ng	100
60) Diethylphthalate	10.298	149	913017	46.824	ng	99
61) 4-Chlorophenyl-phenyle...	10.416	204	422377	44.949	ng	98
62) 4-Nitroaniline	10.445	138	222950	45.463	ng	97
63) Azobenzene	10.575	77	865083	46.068	ng	98
65) 4,6-Dinitro-2-methylph...	10.469	198	155017	49.639	ng	81
66) n-Nitrosodiphenylamine	10.533	169	800889	46.924	ng	99
67) 4-Bromophenyl-phenylether	10.904	248	283319	47.762	ng	# 92
68) Hexachlorobenzene	10.969	284	320333	50.562	ng	95
69) Atrazine	11.063	200	181068	38.991	ng	97
70) Pentachlorophenol	11.163	266	352187	86.559	ng	99
71) Phenanthrene	11.386	178	1318090	46.269	ng	99
72) Anthracene	11.439	178	1354516	46.401	ng	100
73) Carbazole	11.592	167	1177849	46.432	ng	99
74) Di-n-butylphthalate	11.933	149	1362182	46.995	ng	99
75) Fluoranthene	12.580	202	1310021	44.763	ng	99
77) Benzidine	12.704	184	358392	63.380	ng	98
78) Pyrene	12.810	202	1314221	51.380	ng	100
80) Butylbenzylphthalate	13.439	149	466275	50.927	ng	99
81) Benzo(a)anthracene	14.010	228	900609	46.890	ng	99
82) 3,3'-Dichlorobenzidine	13.974	252	288122	47.000	ng	98
83) Chrysene	14.045	228	873113	46.159	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	532049	49.878	ng	98
85) Di-n-octyl phthalate	14.627	149	829334	51.899	ng	100
87) Indeno(1,2,3-cd)pyrene	17.033	276	998694	55.684	ng	99
88) Benzo(b)fluoranthene	15.068	252	816233	50.267	ng	99
89) Benzo(k)fluoranthene	15.104	252	729312	45.393	ng	99
90) Benzo(a)pyrene	15.445	252	698222	46.278	ng	99
91) Dibenzo(a,h)anthracene	17.057	278	823402	56.047	ng	98
92) Benzo(g,h,i)perylene	17.498	276	840738	50.129	ng	98

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

