

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062624\
 Data File : BF138357.D
 Acq On : 26 Jun 2024 17:58
 Operator : CG\JU
 Sample : P2915-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-NORTH-1MSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/27/2024
 Supervised By :mohammad ahmed 06/27/2024

Quant Time: Jun 27 01:54:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 13 01:52:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	213920	20.000	ng	-0.01
21) Naphthalene-d8	8.163	136	834367	20.000	ng	-0.01
39) Acenaphthene-d10	9.916	164	453527	20.000	ng	-0.01
64) Phenanthrene-d10	11.404	188	763863	20.000	ng	-0.01
76) Chrysene-d12	14.039	240	404189	20.000	ng	-0.01
86) Perylene-d12	15.510	264	469969	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1258943	98.448	ng	0.00
7) Phenol-d6	6.516	99	1574986	97.128	ng	0.00
23) Nitrobenzene-d5	7.445	82	989533	69.043	ng	-0.01
42) 2,4,6-Tribromophenol	10.710	330	492305	109.039	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	1885518	66.110	ng	-0.01
79) Terphenyl-d14	12.986	244	1725943	67.632	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.734	88	224167	38.305	ng	96
3) Pyridine	3.487	79	580449	41.732	ng	96
4) n-Nitrosodimethylamine	3.446	42	278943	43.334	ng	90
6) Aniline	6.545	93	561008	35.358	ng	96
8) 2-Chlorophenol	6.669	128	671474	48.568	ng	99
9) Benzaldehyde	6.428	77	128414	14.910	ng	99
10) Phenol	6.528	94	793827	48.239	ng	96
11) bis(2-Chloroethyl)ether	6.616	93	604179	45.874	ng	99
12) 1,3-Dichlorobenzene	6.822	146	691120	43.986	ng	100
13) 1,4-Dichlorobenzene	6.898	146	696345	44.039	ng	99
14) 1,2-Dichlorobenzene	7.051	146	671729	45.254	ng	99
15) Benzyl Alcohol	7.022	79	536829	49.099	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.151	45	757218	40.737	ng	96
17) 2-Methylphenol	7.134	107	518833	47.565	ng	99
18) Hexachloroethane	7.392	117	245407	45.812	ng	95
19) n-Nitroso-di-n-propyla...	7.298	70	417342	43.419	ng	97
20) 3+4-Methylphenols	7.287	107	605375	44.093	ng	97
22) Acetophenone	7.292	105	766580	40.219	ng	# 93
24) Nitrobenzene	7.463	77	645483	43.576	ng	95
25) Isophorone	7.704	82	1179924	45.515	ng	100
26) 2-Nitrophenol	7.781	139	365789	61.411	ng	99
27) 2,4-Dimethylphenol	7.816	122	449436	49.596	ng	98
28) bis(2-Chloroethoxy)met...	7.910	93	738540	45.510	ng	100
29) 2,4-Dichlorophenol	8.022	162	552519	47.357	ng	98
30) 1,2,4-Trichlorobenzene	8.104	180	606591	44.086	ng	98
31) Naphthalene	8.187	128	1806584	43.788	ng	99
32) Benzoic acid	7.945	122	386624	48.407	ng	98
33) 4-Chloroaniline	8.228	127	261002	16.926	ng	99
34) Hexachlorobutadiene	8.298	225	354278	44.026	ng	98
35) Caprolactam	8.610	113	171632m	48.728	ng	
36) 4-Chloro-3-methylphenol	8.716	107	567749	48.875	ng	97
37) 2-Methylnaphthalene	8.875	142	1206522	44.559	ng	98
38) 1-Methylnaphthalene	8.975	142	1123373	42.337	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	557234	42.046	ng	99
41) Hexachlorocyclopentadiene	9.028	237	492924	114.067	ng	98
43) 2,4,6-Trichlorophenol	9.151	196	415335	50.051	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	439749	48.255	ng	98
46) 1,1'-Biphenyl	9.339	154	1399208	42.045	ng	97
47) 2-Chloronaphthalene	9.363	162	1170668	44.511	ng	98
48) 2-Nitroaniline	9.463	65	331133	52.279	ng	88
49) Acenaphthylene	9.781	152	1797137	48.729	ng	99
50) Dimethylphthalate	9.639	163	1397402	50.288	ng	100
51) 2,6-Dinitrotoluene	9.704	165	319724	53.037	ng	87
52) Acenaphthene	9.951	154	1220146m	48.518	ng	
53) 3-Nitroaniline	9.869	138	195341	30.332	ng	96
54) 2,4-Dinitrophenol	9.981	184	307336	94.015	ng	# 88
55) Dibenzofuran	10.128	168	1593293	45.321	ng	98
56) 4-Nitrophenol	10.033	139	470670	92.484	ng	95
57) 2,4-Dinitrotoluene	10.110	165	425662	57.248	ng	94
58) Fluorene	10.469	166	1209356	43.752	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	349782	49.113	ng	99
60) Diethylphthalate	10.339	149	1321200	50.096	ng	99
61) 4-Chlorophenyl-phenyle...	10.457	204	629893	45.756	ng	99
62) 4-Nitroaniline	10.486	138	302019	46.938	ng	97
63) Azobenzene	10.616	77	1193106	45.156	ng	97
65) 4,6-Dinitro-2-methylph...	10.510	198	221759	55.677	ng	93
66) n-Nitrosodiphenylamine	10.580	169	1111406	47.525	ng	100
67) 4-Bromophenyl-phenylether	10.945	248	408066	47.409	ng	# 90
68) Hexachlorobenzene	11.010	284	462752	46.069	ng	99
69) Atrazine	11.104	200	345892	51.928	ng	99
70) Pentachlorophenol	11.204	266	488309	83.354	ng	97
71) Phenanthrene	11.427	178	1725867	45.687	ng	99
72) Anthracene	11.480	178	1753255	46.746	ng	99
73) Carbazole	11.633	167	1483445	43.123	ng	99
74) Di-n-butylphthalate	11.963	149	1913964	53.744	ng	99
75) Fluoranthene	12.616	202	1594696	41.398	ng	98
77) Benzidine	12.733	184	468212	100.149	ng	98
78) Pyrene	12.845	202	1595490	45.764	ng	98
80) Butylbenzylphthalate	13.457	149	623833	65.783	ng	97
81) Benzo(a)anthracene	14.027	228	1255430	48.415	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	313429	43.893	ng	98
83) Chrysene	14.068	228	1218284	49.290	ng	100
84) Bis(2-ethylhexyl)phtha...	14.016	149	808452	73.299	ng	99
85) Di-n-octyl phthalate	14.627	149	1352017	82.513	ng	97
87) Indeno(1,2,3-cd)pyrene	16.986	276	1414898	46.182	ng	100
88) Benzo(b)fluoranthene	15.080	252	1318911	47.589	ng	100
89) Benzo(k)fluoranthene	15.110	252	1278356	47.066	ng	99
90) Benzo(a)pyrene	15.451	252	1224658	51.436	ng	99
91) Dibenzo(a,h)anthracene	17.004	278	1159196	46.096	ng	99
92) Benzo(g,h,i)perylene	17.433	276	1042537	40.074	ng	99

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

