

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
 Data File : BF136913.D
 Acq On : 03 Jan 2024 18:09
 Operator : CG\JU
 Sample : PB157954BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB157954BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 01/04/2024
 Supervised By :mohammad ahmed 01/04/2024

Quant Time: Jan 04 03:39:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	77590	20.000	ng	-0.01
21) Naphthalene-d8	8.057	136	304111	20.000	ng	-0.01
39) Acenaphthene-d10	9.816	164	158162	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	307337	20.000	ng	0.00
76) Chrysene-d12	13.963	240	165828	20.000	ng	0.00
86) Perylene-d12	15.445	264	141574	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	560707	112.755	ng	0.00
7) Phenol-d6	6.440	99	701607	108.885	ng	0.00
23) Nitrobenzene-d5	7.351	82	429360	72.246	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	215552	115.720	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	870679	74.041	ng	-0.01
79) Terphenyl-d14	12.904	244	962577	81.118	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.616	88	71066	34.299	ng	96
3) Pyridine	3.381	79	163795	32.400	ng	94
4) n-Nitrosodimethylamine	3.346	42	96647	35.800	ng	99
6) Aniline	6.451	93	225359	34.981	ng	# 34
8) 2-Chlorophenol	6.575	128	216573	39.797	ng	94
9) Benzaldehyde	6.340	77	91787	25.044	ng	99
10) Phenol	6.451	94	264075	38.391	ng	88
11) bis(2-Chloroethyl)ether	6.522	93	205722	39.327	ng	97
12) 1,3-Dichlorobenzene	6.722	146	218995	36.889	ng	97
13) 1,4-Dichlorobenzene	6.798	146	223995	36.928	ng	98
14) 1,2-Dichlorobenzene	6.951	146	213552	37.835	ng	97
15) Benzyl Alcohol	6.934	79	172990	39.795	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.057	45	318804	37.277	ng	87
17) 2-Methylphenol	7.045	107	171969	38.145	ng	99
18) Hexachloroethane	7.287	117	78820	35.780	ng	93
19) n-Nitroso-di-n-propyla...	7.198	70	146679	37.552	ng	96
20) 3+4-Methylphenols	7.198	107	217493	38.616	ng	# 83
22) Acetophenone	7.192	105	296219	38.861	ng	# 90
24) Nitrobenzene	7.369	77	215113	36.133	ng	93
25) Isophorone	7.604	82	404329	37.390	ng	100
26) 2-Nitrophenol	7.681	139	114534	40.211	ng	99
27) 2,4-Dimethylphenol	7.722	122	190438	54.918	ng	98
28) bis(2-Chloroethoxy)met...	7.816	93	252377	38.394	ng	100
29) 2,4-Dichlorophenol	7.928	162	177615	39.785	ng	96
30) 1,2,4-Trichlorobenzene	8.004	180	185633	37.320	ng	99
31) Naphthalene	8.081	128	626201	37.481	ng	99
32) Benzoic acid	7.875	122	131175	39.096	ng	97
33) 4-Chloroaniline	8.134	127	130341	21.130	ng	99
34) Hexachlorobutadiene	8.192	225	107769	35.660	ng	99
35) Caprolactam	8.516	113	61540m	41.785	ng	
36) 4-Chloro-3-methylphenol	8.633	107	195720	38.942	ng	98
37) 2-Methylnaphthalene	8.775	142	406400	37.797	ng	100
38) 1-Methylnaphthalene	8.875	142	375806	35.626	ng	100
40) 1,2,4,5-Tetrachloroben...	8.939	216	194497	39.708	ng	99
41) Hexachlorocyclopentadiene	8.922	237	143085	90.906	ng	99
43) 2,4,6-Trichlorophenol	9.057	196	126068	40.132	ng	99

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44) 2,4,5-Trichlorophenol	9.110	196	136438	38.976	ng	97
46) 1,1'-Biphenyl	9.239	154	521084	39.261	ng	100
47) 2-Chloronaphthalene	9.263	162	381042	37.765	ng	99
48) 2-Nitroaniline	9.369	65	122478	38.764	ng	94
49) Acenaphthylene	9.681	152	607357	39.382	ng	100
50) Dimethylphthalate	9.545	163	454584	39.349	ng	100
51) 2,6-Dinitrotoluene	9.610	165	101323	38.645	ng	# 87
52) Acenaphthene	9.851	154	369670	38.669	ng	98
53) 3-Nitroaniline	9.780	138	80825	29.101	ng	94
54) 2,4-Dinitrophenol	9.898	184	99345	68.911	ng	93
55) Dibenzofuran	10.028	168	556113	38.375	ng	97
56) 4-Nitrophenol	9.963	139	130716	69.719	ng	96
57) 2,4-Dinitrotoluene	10.022	165	126929	37.292	ng	95
58) Fluorene	10.369	166	454121	39.494	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.151	232	114739	42.580	ng	98
60) Diethylphthalate	10.245	149	465491	38.835	ng	99
61) 4-Chlorophenyl-phenyle...	10.363	204	216612	38.757	ng	93
62) 4-Nitroaniline	10.404	138	100349	37.479	ng	96
63) Azobenzene	10.522	77	427273	38.165	ng	99
65) 4,6-Dinitro-2-methylph...	10.433	198	73224	41.573	ng	85
66) n-Nitrosodiphenylamine	10.486	169	403563	40.422	ng	99
67) 4-Bromophenyl-phenylether	10.851	248	132161	38.720	ng	99
68) Hexachlorobenzene	10.916	284	148671	39.086	ng	96
69) Atrazine	11.022	200	181482	71.071	ng	98
70) Pentachlorophenol	11.122	266	116838	65.900	ng	98
71) Phenanthrene	11.339	178	640170	40.595	ng	99
72) Anthracene	11.386	178	658953	41.261	ng	99
73) Carbazole	11.551	167	608853	40.436	ng	100
74) Di-n-butylphthalate	11.874	149	744579	40.525	ng	100
75) Fluoranthene	12.527	202	667437	39.568	ng	97
77) Benzidine	12.657	184	149935	30.664	ng	99
78) Pyrene	12.763	202	668575	41.070	ng	99
80) Butylbenzylphthalate	13.380	149	255039	43.056	ng	97
81) Benzo(a)anthracene	13.957	228	468557	40.691	ng	99
82) 3,3'-Dichlorobenzidine	13.921	252	126903	38.807	ng	97
83) Chrysene	13.992	228	442749	39.319	ng	98
84) Bis(2-ethylhexyl)phtha...	13.945	149	338312	45.394	ng	# 98
85) Di-n-octyl phthalate	14.557	149	482679	41.869	ng	97
87) Indeno(1,2,3-cd)pyrene	16.951	276	476980	46.663	ng	100
88) Benzo(b)fluoranthene	15.010	252	399712	42.271	ng	98
89) Benzo(k)fluoranthene	15.039	252	357873	40.053	ng	99
90) Benzo(a)pyrene	15.380	252	331751	41.564	ng	98
91) Dibenzo(a,h)anthracene	16.968	278	397555	47.407	ng	98
92) Benzo(g,h,i)perylene	17.409	276	405615	47.225	ng	99

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

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