

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
 Data File : BF140648.D
 Acq On : 26 Nov 2024 18:00
 Operator : RC/JU
 Sample : P4938-04MSD
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-732MSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Nov 27 00:07:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 15:23:48 2024
 Response via : Initial Calibration

11/27/2024
 Supervised By :mohammad
 ahmed

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	65801	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	240817	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	128112	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	225568	20.000	ng	0.00
76) Chrysene-d12	14.051	240	155066	20.000	ng	0.00
86) Perylene-d12	15.545	264	144332	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	560561	145.353	ng	0.01
7) Phenol-d6	6.516	99	702751	137.836	ng	0.00
23) Nitrobenzene-d5	7.434	82	478711	101.680	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	202121	147.516	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	863754	100.455	ng	0.00
79) Terphenyl-d14	12.980	244	804394	80.775	ng	-0.01

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.757	88	65999	40.703	ng	# 83
3) Pyridine	3.504	79	108451	30.399	ng	97
4) n-Nitrosodimethylamine	3.422	42	95192	45.398	ng	92
6) Aniline	6.539	93	49507	13.796	ng	# 50
8) 2-Chlorophenol	6.663	128	209526	50.499	ng	95
9) Benzaldehyde	6.428	77	16928	6.272	ng	97
10) Phenol	6.528	94	247754	47.583	ng	85
11) bis(2-Chloroethyl)ether	6.604	93	193752	48.628	ng	100
12) 1,3-Dichlorobenzene	6.810	146	194828	41.790	ng	98
13) 1,4-Dichlorobenzene	6.892	146	201220	42.626	ng	99
14) 1,2-Dichlorobenzene	7.039	146	192045	43.416	ng	99
15) Benzyl Alcohol	7.016	79	174263	46.089	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.139	45	240250	51.040	ng	88
17) 2-Methylphenol	7.134	107	169206	50.946	ng	98
18) Hexachloroethane	7.381	117	71036	40.291	ng	97
19) n-Nitroso-di-n-propyla...	7.281	70	148864	49.404	ng	98
20) 3+4-Methylphenols	7.286	107	223859	52.438	ng	91
22) Acetophenone	7.281	105	292973	49.902	ng	96
24) Nitrobenzene	7.457	77	225745	46.393	ng	97
25) Isophorone	7.692	82	396618	50.508	ng	100
26) 2-Nitrophenol	7.769	139	110078	51.048	ng	97
27) 2,4-Dimethylphenol	7.810	122	171501m	66.473	ng	
28) bis(2-Chloroethoxy)met...	7.898	93	236410	49.418	ng	99
29) 2,4-Dichlorophenol	8.016	162	181404	53.062	ng	97
30) 1,2,4-Trichlorobenzene	8.092	180	177279	45.416	ng	98
31) Naphthalene	8.175	128	602153	48.544	ng	99
32) Benzoic acid	7.933	122	116187	51.188	ng	99
33) 4-Chloroaniline	8.233	127	34998	9.424	ng	97
34) Hexachlorobutadiene	8.286	225	111029	42.944	ng	98
35) Caprolactam	8.586	113	49775m	47.047	ng	
36) 4-Chloro-3-methylphenol	8.716	107	199297	52.067	ng	98
37) 2-Methylnaphthalene	8.863	142	397926	50.507	ng	100
38) 1-Methylnaphthalene	8.963	142	367223	47.552	ng	100
40) 1,2,4,5-Tetrachloroben...	9.033	216	192792	51.405	ng	99
41) Hexachlorocyclopentadiene	9.016	237	94787	109.480	ng	98
43) 2,4,6-Trichlorophenol	9.145	196	128858	54.754	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	138110	54.074	ng	95
46) 1,1'-Biphenyl	9.328	154	502372	52.522	ng	99
47) 2-Chloronaphthalene	9.357	162	372782	51.435	ng	98
48) 2-Nitroaniline	9.451	65	122343	52.591	ng	98
49) Acenaphthylene	9.769	152	609254	55.637	ng	100
50) Dimethylphthalate	9.627	163	458727	54.368	ng	100
51) 2,6-Dinitrotoluene	9.692	165	97739	51.077	ng	95
52) Acenaphthene	9.939	154	368259	52.927	ng	99
53) 3-Nitroaniline	9.869	138	20696	11.058	ng	98
54) 2,4-Dinitrophenol	9.980	184	94162	93.545	ng	94
55) Dibenzofuran	10.116	168	563810	53.124	ng	99
56) 4-Nitrophenol	10.039	139	132292	103.646	ng	93
57) 2,4-Dinitrotoluene	10.104	165	130813	51.437	ng	98
58) Fluorene	10.457	166	441119	51.782	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	116842	59.525	ng	93
60) Diethylphthalate	10.327	149	446963	52.139	ng	100
61) 4-Chlorophenyl-phenyle...	10.445	204	217783	52.093	ng	98
62) 4-Nitroaniline	10.480	138	69362	35.336	ng	93
63) Azobenzene	10.610	77	400385	49.320	ng	97
65) 4,6-Dinitro-2-methylph...	10.510	198	65975	57.237	ng	97
66) n-Nitrosodiphenylamine	10.569	169	311765	46.737	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	125100	53.853	ng	97
68) Hexachlorobenzene	11.010	284	137408	51.085	ng	99
69) Atrazine	11.092	200	73140	38.975	ng	99
70) Pentachlorophenol	11.210	266	138877	117.310	ng	97
71) Phenanthrene	11.421	178	592030	54.601	ng	99
72) Anthracene	11.474	178	600835	56.649	ng	100
73) Carbazole	11.633	167	511892	50.169	ng	99
74) Di-n-butylphthalate	11.951	149	686093	58.243	ng	100
75) Fluoranthene	12.616	202	609949	51.811	ng	100
77) Benzidine	12.745	184	58487	12.973	ng	99
78) Pyrene	12.845	202	616429	42.993	ng	99
80) Butylbenzylphthalate	13.457	149	247037	47.829	ng	99
81) Benzo(a)anthracene	14.039	228	563828	54.929	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	28322	9.190	ng	100
83) Chrysene	14.074	228	509775	54.312	ng	100
84) Bis(2-ethylhexyl)phtha...	14.015	149	332264	50.946	ng	99
85) Di-n-octyl phthalate	14.639	149	546234	61.285	ng	98
87) Indeno(1,2,3-cd)pyrene	17.051	276	386990	41.143	ng	97
88) Benzo(b)fluoranthene	15.109	252	502250	55.401	ng	99
89) Benzo(k)fluoranthene	15.139	252	468484	59.053	ng	100
90) Benzo(a)pyrene	15.480	252	428040	58.070	ng	98
91) Dibenzo(a,h)anthracene	17.062	278	325450	42.249	ng	99
92) Benzo(g,h,i)perylene	17.509	276	282466	35.934	ng	98

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

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