

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091624\
 Data File : BF139385.D
 Acq On : 16 Sep 2024 18:23
 Operator : RC/JU
 Sample : P3947-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RW-1MSD

Quant Time: Sep 17 01:37:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	58657	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	210860	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	103397	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	155991	20.000	ng	0.00
76) Chrysene-d12	14.033	240	92472	20.000	ng	0.00
86) Perylene-d12	15.498	264	73636	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	423232	117.708	ng	0.01
7) Phenol-d6	6.510	99	550215	116.584	ng	0.00
23) Nitrobenzene-d5	7.445	82	380087	84.741	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	113124	102.767	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	618944	83.956	ng	0.00
79) Terphenyl-d14	12.980	244	417283	74.068	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.740	88	58290	40.407	ng	100
3) Pyridine	3.481	79	146330	46.008	ng	96
4) n-Nitrosodimethylamine	3.440	42	130780	47.440	ng	100
6) Aniline	6.545	93	123596	34.702	ng	96
8) 2-Chlorophenol	6.663	128	178073	48.112	ng	99
9) Benzaldehyde	6.428	77	38640	14.171	ng	98
10) Phenol	6.528	94	231217	48.103	ng	95
11) bis(2-Chloroethyl)ether	6.616	93	174522	47.770	ng	99
12) 1,3-Dichlorobenzene	6.822	146	194580	46.226	ng	100
13) 1,4-Dichlorobenzene	6.898	146	195749	46.019	ng	100
14) 1,2-Dichlorobenzene	7.045	146	184292	45.851	ng	99
15) Benzyl Alcohol	7.022	79	183654	48.775	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.151	45	303221	53.855	ng	97
17) 2-Methylphenol	7.134	107	142163	48.344	ng	100
18) Hexachloroethane	7.392	117	74468	45.963	ng	99
19) n-Nitroso-di-n-propyla...	7.292	70	145290	51.950	ng	99
20) 3+4-Methylphenols	7.286	107	193379	48.200	ng	# 89
22) Acetophenone	7.286	105	269087	48.814	ng	97
24) Nitrobenzene	7.463	77	216224	46.839	ng	99
25) Isophorone	7.698	82	363754	51.280	ng	98
26) 2-Nitrophenol	7.775	139	91764	51.023	ng	98
27) 2,4-Dimethylphenol	7.816	122	143914	60.662	ng	99
28) bis(2-Chloroethoxy)met...	7.910	93	209298	51.771	ng	98
29) 2,4-Dichlorophenol	8.016	162	144531	47.956	ng	99
30) 1,2,4-Trichlorobenzene	8.098	180	161809	46.078	ng	98
31) Naphthalene	8.186	128	527657	47.933	ng	100
32) Benzoic acid	7.934	122	100154	45.112	ng	99
33) 4-Chloroaniline	8.233	127	38923	11.558	ng	99
34) Hexachlorobutadiene	8.298	225	112329	47.083	ng	99
35) Caprolactam	8.598	113	42262	45.586	ng	96
36) 4-Chloro-3-methylphenol	8.710	107	158129	45.909	ng	99
37) 2-Methylnaphthalene	8.875	142	335398	48.002	ng	100
38) 1-Methylnaphthalene	8.975	142	307627	44.845	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	160144	52.260	ng	99
41) Hexachlorocyclopentadiene	9.022	237	111650	104.351	ng	99
43) 2,4,6-Trichlorophenol	9.151	196	102312	51.319	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	102439	48.711	ng	99
46) 1,1'-Biphenyl	9.339	154	411676	51.753	ng	100
47) 2-Chloronaphthalene	9.363	162	294372	49.494	ng	98
48) 2-Nitroaniline	9.457	65	109312	49.413	ng	99
49) Acenaphthylene	9.775	152	468990	52.384	ng	100
50) Dimethylphthalate	9.639	163	360922	53.183	ng	100
51) 2,6-Dinitrotoluene	9.698	165	72055	47.928	ng	99
52) Acenaphthene	9.945	154	317875	50.872	ng	99
53) 3-Nitroaniline	9.863	138	47680	31.734	ng	95
54) 2,4-Dinitrophenol	9.969	184	54408	64.947	ng	97
55) Dibenzofuran	10.122	168	416015	47.767	ng	97
56) 4-Nitrophenol	10.027	139	99269	79.848	ng	93
57) 2,4-Dinitrotoluene	10.098	165	91488	45.075	ng	98
58) Fluorene	10.463	166	325108	45.751	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.233	232	77669	44.647	ng	98
60) Diethylphthalate	10.333	149	349864	49.957	ng	99
61) 4-Chlorophenyl-phenyle...	10.451	204	161419	46.742	ng	100
62) 4-Nitroaniline	10.475	138	58383	35.604	ng	99
63) Azobenzene	10.610	77	357744	51.936	ng	98
65) 4,6-Dinitro-2-methylph...	10.504	198	33637	40.856	ng	98
66) n-Nitrosodiphenylamine	10.569	169	264084	57.905	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	86930	56.235	ng	100
68) Hexachlorobenzene	11.004	284	92925	51.220	ng	98
69) Atrazine	11.098	200	80471	74.116	ng	98
70) Pentachlorophenol	11.198	266	109192	100.304	ng	98
71) Phenanthrene	11.422	178	489574	65.052	ng	99
72) Anthracene	11.474	178	409426	55.642	ng	99
73) Carbazole	11.627	167	349493	49.626	ng	99
74) Di-n-butylphthalate	11.957	149	443766	56.565	ng	100
75) Fluoranthene	12.610	202	528350	65.739	ng	100
77) Benzidine	12.727	184	79648	60.118	ng	99
78) Pyrene	12.839	202	527119	66.388	ng	100
80) Butylbenzylphthalate	13.451	149	162905	60.277	ng	99
81) Benzo(a)anthracene	14.021	228	363951	59.417	ng	99
82) 3,3'-Dichlorobenzidine	13.986	252	65011	36.242	ng	99
83) Chrysene	14.063	228	331092	58.753	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	209390	60.903	ng	99
85) Di-n-octyl phthalate	14.621	149	305395	60.316	ng	99
87) Indeno(1,2,3-cd)pyrene	16.968	276	207031	47.282	ng	98
88) Benzo(b)fluoranthene	15.068	252	274928	60.538	ng	99
89) Benzo(k)fluoranthene	15.098	252	224182	53.539	ng	100
90) Benzo(a)pyrene	15.439	252	223698	59.393	ng	99
91) Dibenzo(a,h)anthracene	16.986	278	161747	44.704	ng	99
92) Benzo(g,h,i)perylene	17.409	276	155354	43.365	ng	98

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

