

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042524\
 Data File : BG061040.D
 Acq On : 25 Apr 2024 16:40
 Operator : MA/JU
 Sample : P2261-01MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RO-RBR200059MS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/26/2024
 Supervised By :mohammad ahmed 04/27/2024

Quant Time: Apr 25 17:39:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 22 04:22:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	28067	20.000	ng	0.00
21) Naphthalene-d8	10.738	136	107691	20.000	ng	0.00
39) Acenaphthene-d10	14.574	164	81719	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	188224	20.000	ng	0.00
76) Chrysene-d12	21.584	240	189029	20.000	ng	0.00
86) Perylene-d12	24.715	264	173721	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.538	112	182381	131.439	ng	0.00
7) Phenol-d6	7.118	99	256560	124.618	ng	0.00
23) Nitrobenzene-d5	9.098	82	157508	65.713	ng	0.00
42) 2,4,6-Tribromophenol	16.067	330	194059	108.142	ng	0.00
45) 2-Fluorobiphenyl	13.194	172	420531	71.222	ng	0.00
79) Terphenyl-d14	19.939	244	719422	59.076	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.452	88	18552	38.632	ng	93
3) Pyridine	3.840	79	52751	35.537	ng	99
4) n-Nitrosodimethylamine	3.758	42	60426	45.261	ng	99
6) Aniline	7.271	93	44898	18.145	ng	98
8) 2-Chlorophenol	7.512	128	89065	51.515	ng	96
9) Benzaldehyde	7.083	77	3031	2.862	ng	# 80
10) Phenol	7.148	94	105201	51.432	ng	96
11) bis(2-Chloroethyl)ether	7.365	93	70082	47.889	ng	93
12) 1,3-Dichlorobenzene	7.835	146	87091	45.424	ng	99
13) 1,4-Dichlorobenzene	7.976	146	88307	44.715	ng	97
14) 1,2-Dichlorobenzene	8.294	146	86649	44.895	ng	96
15) Benzyl Alcohol	8.182	79	74622	38.389	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.464	45	126172	39.606	ng	99
17) 2-Methylphenol	8.388	107	62193	37.657	ng	91
18) Hexachloroethane	9.022	117	31813	44.706	ng	95
19) n-Nitroso-di-n-propyla...	8.746	70	55926	36.451	ng	98
20) 3+4-Methylphenols	8.717	107	88274	38.217	ng	97
22) Acetophenone	8.764	105	121981	40.564	ng	# 93
24) Nitrobenzene	9.140	77	90893	39.463	ng	95
25) Isophorone	9.662	82	154130	38.424	ng	98
26) 2-Nitrophenol	9.851	139	43340	40.037	ng	98
27) 2,4-Dimethylphenol	9.915	122	58845	33.915	ng	94
28) bis(2-Chloroethoxy)met...	10.144	93	83027	39.346	ng	97
29) 2,4-Dichlorophenol	10.391	162	86102	42.997	ng	99
30) 1,2,4-Trichlorobenzene	10.603	180	103654	46.914	ng	96
31) Naphthalene	10.791	128	258878	43.909	ng	99
32) Benzoic acid	10.068	122	53529m	40.491	ng	
33) 4-Chloroaniline	10.890	127	39525	14.758	ng	99
34) Hexachlorobutadiene	11.079	225	90548	46.600	ng	97
35) Caprolactam	11.678	113	23990m	37.612	ng	
36) 4-Chloro-3-methylphenol	12.030	107	93682	40.112	ng	95
37) 2-Methylnaphthalene	12.400	142	184468	39.794	ng	99
38) 1-Methylnaphthalene	12.618	142	173962	40.215	ng	98
40) 1,2,4,5-Tetrachloroben...	12.765	216	150969	51.985	ng	99
41) Hexachlorocyclopentadiene	12.747	237	61001	35.699	ng	96
43) 2,4,6-Trichlorophenol	13.006	196	92041	48.214	ng	96

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44) 2,4,5-Trichlorophenol	13.088	196	101342	44.878	ng	97
46) 1,1'-Biphenyl	13.405	154	262723	47.729	ng	100
47) 2-Chloronaphthalene	13.446	162	207123	49.076	ng	97
48) 2-Nitroaniline	13.646	65	74826	43.929	ng	93
49) Acenaphthylene	14.298	152	340675	47.929	ng	99
50) Dimethylphthalate	14.034	163	281391	41.758	ng	97
51) 2,6-Dinitrotoluene	14.145	165	57482	40.882	ng	98
52) Acenaphthene	14.639	154	187639	43.445	ng	95
53) 3-Nitroaniline	14.475	138	41848	30.184	ng	99
54) 2,4-Dinitrophenol	14.686	184	24183	30.128	ng	92
55) Dibenzofuran	14.974	168	324216	43.472	ng	98
56) 4-Nitrophenol	14.798	139	87148	87.738	ng	95
57) 2,4-Dinitrotoluene	14.939	165	82925	40.949	ng	93
58) Fluorene	15.620	166	270192	42.922	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.203	232	100727	44.196	ng	# 98
60) Diethylphthalate	15.397	149	268623	40.478	ng	98
61) 4-Chlorophenyl-phenyle...	15.614	204	160527	41.347	ng	99
62) 4-Nitroaniline	15.644	138	55941	37.419	ng	96
63) Azobenzene	15.908	77	222028	41.133	ng	98
65) 4,6-Dinitro-2-methylph...	15.702	198	19316	16.620	ng	96
66) n-Nitrosodiphenylamine	15.832	169	241631	46.596	ng	99
67) 4-Bromophenyl-phenylether	16.513	248	114729	46.185	ng	99
68) Hexachlorobenzene	16.631	284	137149	49.984	ng	98
69) Atrazine	16.784	200	101369	52.230	ng	95
70) Pentachlorophenol	16.977	266	170579	94.807	ng	98
71) Phenanthrene	17.365	178	444436	46.259	ng	99
72) Anthracene	17.453	178	442413	44.711	ng	98
73) Carbazole	17.724	167	366655	44.617	ng	99
74) Di-n-butylphthalate	18.288	149	413456	41.440	ng	100
75) Fluoranthene	19.381	202	630689	49.464	ng	97
77) Benzidine	19.557	184	222029	63.392	ng	99
78) Pyrene	19.739	202	662001	51.434	ng	99
80) Butylbenzylphthalate	20.632	149	175667	40.948	ng	97
81) Benzo(a)anthracene	21.566	228	657345	49.462	ng	96
82) 3,3'-Dichlorobenzidine	21.478	252	174282	36.287	ng	100
83) Chrysene	21.631	228	636227	50.283	ng	97
84) Bis(2-ethylhexyl)phtha...	21.490	149	259293	41.573	ng	99
85) Di-n-octyl phthalate	22.671	149	438793	42.375	ng	98
87) Indeno(1,2,3-cd)pyrene	28.264	276	463917m	38.496	ng	
88) Benzo(b)fluoranthene	23.728	252	542120	55.067	ng	99
89) Benzo(k)fluoranthene	23.793	252	608446	61.224	ng	97
90) Benzo(a)pyrene	24.569	252	538505	56.513	ng	96
91) Dibenzo(a,h)anthracene	28.329	278	356113	36.169	ng	99
92) Benzo(g,h,i)perylene	29.375	276	302968m	31.034	ng	

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

