

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
 Data File : BG056011.D
 Acq On : 19 Dec 2022 13:03
 Operator : CG/JU
 Sample : PB149627BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB149627BS

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 12/20/2022
 Supervised By :mohammad ahmed 12/20/2022

Quant Time: Dec 20 01:05:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 20 01:01:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.371	152	7627	20.000	ng	0.00
21) Naphthalene-d8	11.203	136	32437	20.000	ng	# 0.00
39) Acenaphthene-d10	14.975	164	30483	20.000	ng	0.00
64) Phenanthrene-d10	17.713	188	91639	20.000	ng	0.00
76) Chrysene-d12	22.025	240	102427	20.000	ng	0.00
86) Perylene-d12	25.515	264	108541	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.885	112	48872	150.273	ng	0.00
7) Phenol-d6	7.495	99	60957	145.733	ng	0.00
23) Nitrobenzene-d5	9.540	82	39801	85.747	ng	0.00
42) 2,4,6-Tribromophenol	16.455	330	97023	131.998	ng	0.00
45) 2-Fluorobiphenyl	13.606	172	153401	70.578	ng	0.00
79) Terphenyl-d14	20.286	244	452319	78.048	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.711	88	3590	37.931	ng	98
3) Pyridine	4.123	79	9799	31.460	ng	94
4) n-Nitrosodimethylamine	4.034	42	10256	36.629	ng	# 93
6) Aniline	7.677	93	18195	36.332	ng	93
8) 2-Chlorophenol	7.924	128	20272	48.315	ng	97
9) Benzaldehyde	7.495	77	4859	18.497	ng	99
10) Phenol	7.524	94	22291	48.693	ng	96
11) bis(2-Chloroethyl)ether	7.765	93	12958	43.456	ng	93
12) 1,3-Dichlorobenzene	8.259	146	22377	41.228	ng	93
13) 1,4-Dichlorobenzene	8.406	146	23174	41.779	ng	95
14) 1,2-Dichlorobenzene	8.729	146	22329	42.122	ng	96
15) Benzyl Alcohol	8.594	79	19074	43.886	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.888	45	17987	46.347	ng	94
17) 2-Methylphenol	8.794	107	17047	47.329	ng	94
18) Hexachloroethane	9.463	117	7395	43.920	ng	97
19) n-Nitroso-di-n-propyla...	9.170	70	12681	42.296	ng	95
20) 3+4-Methylphenols	9.123	107	23030	44.806	ng	94
22) Acetophenone	9.193	105	30615	38.074	ng	96
24) Nitrobenzene	9.587	77	20564	44.014	ng	95
25) Isophorone	10.104	82	37723	38.757	ng	99
26) 2-Nitrophenol	10.298	139	12158	43.958	ng	97
27) 2,4-Dimethylphenol	10.339	122	21528m	46.347	ng	
28) bis(2-Chloroethoxy)met...	10.586	93	19129	44.469	ng	97
29) 2,4-Dichlorophenol	10.832	162	26853	40.730	ng	94
30) 1,2,4-Trichlorobenzene	11.056	180	31181	36.383	ng	98
31) Naphthalene	11.255	128	62680	36.622	ng	98
32) Benzoic acid	10.450	122	16293	41.732	ng	90
33) 4-Chloroaniline	11.344	127	18340	24.996	ng	98
34) Hexachlorobutadiene	11.532	225	28249	35.060	ng	94
35) Caprolactam	12.113	113	7295m	39.714	ng	
36) 4-Chloro-3-methylphenol	12.436	107	25490	41.232	ng	92
37) 2-Methylnaphthalene	12.830	142	54267	35.888	ng	96
38) 1-Methylnaphthalene	13.047	142	51401	34.893	ng	98
40) 1,2,4,5-Tetrachloroben...	13.188	216	48502	37.712	ng	99
41) Hexachlorocyclopentadiene	13.171	237	67805	92.930	ng	99
43) 2,4,6-Trichlorophenol	13.418	196	31203	40.985	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.488	196	35015	40.800	ng	98
46) 1,1'-Biphenyl	13.817	154	80668	39.042	ng	98
47) 2-Chloronaphthalene	13.864	162	63942	38.086	ng	95
48) 2-Nitroaniline	14.052	65	15845	47.883	ng	92
49) Acenaphthylene	14.704	152	101227	37.774	ng	97
50) Dimethylphthalate	14.416	163	98157	37.245	ng	99
51) 2,6-Dinitrotoluene	14.534	165	19600	48.059	ng	92
52) Acenaphthene	15.045	154	79003	43.914	ng	99
53) 3-Nitroaniline	14.869	138	13360	32.206	ng	# 89
54) 2,4-Dinitrophenol	15.069	184	28912	95.360	ng	# 79
55) Dibenzofuran	15.374	168	114041	37.696	ng	96
56) 4-Nitrophenol	15.157	139	31803	95.671	ng	98
57) 2,4-Dinitrotoluene	15.315	165	29493	47.909	ng	# 84
58) Fluorene	16.015	166	93212	37.691	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.586	232	40130	40.930	ng	99
60) Diethylphthalate	15.768	149	95381	37.028	ng	97
61) 4-Chlorophenyl-phenyle...	15.997	204	63538	37.714	ng	99
62) 4-Nitroaniline	16.026	138	19531	43.234	ng	97
63) Azobenzene	16.291	77	59671	38.060	ng	96
65) 4,6-Dinitro-2-methylph...	16.079	198	18830	43.014	ng	96
66) n-Nitrosodiphenylamine	16.208	169	89194	37.142	ng	98
67) 4-Bromophenyl-phenylether	16.896	248	48387	36.034	ng	94
68) Hexachlorobenzene	17.019	284	60180	36.385	ng	96
69) Atrazine	17.143	200	38365	35.283	ng	97
70) Pentachlorophenol	17.354	266	80521	85.156	ng	93
71) Phenanthrene	17.760	178	176724	36.799	ng	99
72) Anthracene	17.848	178	177158	37.288	ng	99
73) Carbazole	18.106	167	149933	38.263	ng	99
74) Di-n-butylphthalate	18.641	149	165445	36.306	ng	99
75) Fluoranthene	19.745	202	239991	35.688	ng	99
77) Benzidine	19.910	184	80787	44.912	ng	99
78) Pyrene	20.104	202	242783	39.172	ng	99
80) Butylbenzylphthalate	20.973	149	70218	39.981	ng	98
81) Benzo(a)anthracene	22.002	228	264489	36.696	ng	98
82) 3,3'-Dichlorobenzidine	21.902	252	90837	35.029	ng	98
83) Chrysene	22.072	228	247611	36.656	ng	99
84) Bis(2-ethylhexyl)phtha...	21.878	149	98892	37.900	ng	99
85) Di-n-octyl phthalate	23.188	149	168496	38.069	ng	100
87) Indeno(1,2,3-cd)pyrene	29.546	276	330895	37.911	ng	99
88) Benzo(b)fluoranthene	24.399	252	296548	42.215	ng	98
89) Benzo(k)fluoranthene	24.475	252	281131	40.112	ng	97
90) Benzo(a)pyrene	25.351	252	255198	42.979	ng	# 99
91) Dibenzo(a,h)anthracene	29.616	278	288433	39.482	ng	# 96
92) Benzo(g,h,i)perylene	30.809	276	278132	39.377	ng	96

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

