

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072723\
 Data File : BP016394.D
 Acq On : 27 Jul 2023 11:37
 Operator : MA/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/28/2023
 Supervised By :mohammad ahmed 07/28/2023

Quant Time: Jul 27 17:35:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 27 17:32:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	380607	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	1613288	20.000	ng	0.00
39) Acenaphthene-d10	14.734	164	1014199	20.000	ng	0.00
64) Phenanthrene-d10	17.498	188	2182918	20.000	ng	0.00
76) Chrysene-d12	21.586	240	1991981	20.000	ng	# 0.00
86) Perylene-d12	24.186	264	2209674	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.599	112	241254	10.341	ng	0.00
7) Phenol-d6	7.222	99	328494	10.273	ng	0.00
23) Nitrobenzene-d5	9.287	82	278170	9.652	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0d	0.000	ng	
45) 2-Fluorobiphenyl	13.357	172	742829	10.501	ng	0.00
79) Terphenyl-d14	20.039	244	1049845	11.133	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.464	88	49241	5.412	ng	94
3) Pyridine	3.875	79	143753	5.370	ng	99
4) n-Nitrosodimethylamine	3.805	42	54922	5.356	ng	# 97
6) Aniline	7.416	93	203960	5.164	ng	96
8) 2-Chlorophenol	7.640	128	125576	5.134	ng	96
9) Benzaldehyde	7.240	77	110905	3.740	ng	94
10) Phenol	7.246	94	161333	5.195	ng	98
11) bis(2-Chloroethyl)ether	7.522	93	136635	5.413	ng	99
12) 1,3-Dichlorobenzene	7.975	146	139603	5.316	ng	95
13) 1,4-Dichlorobenzene	8.128	146	142084	5.334	ng	98
14) 1,2-Dichlorobenzene	8.446	146	136799	5.334	ng	98
15) Benzyl Alcohol	8.340	79	109250	5.001	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.628	45	160526	5.607	ng	98
17) 2-Methylphenol	8.522	107	114366	5.110	ng	99
18) Hexachloroethane	9.169	117	50417	5.306	ng	95
19) n-Nitroso-di-n-propyla...	8.904	70	105976	5.290	ng	95
20) 3+4-Methylphenols	8.851	107	152623	5.021	ng	98
22) Acetophenone	8.934	105	210151	5.159	ng	# 97
24) Nitrobenzene	9.328	77	147501	4.983	ng	94
25) Isophorone	9.840	82	289341	4.973	ng	99
26) 2-Nitrophenol	10.034	139	41534	6.169	ng	# 86
27) 2,4-Dimethylphenol	10.069	122	128406	4.933	ng	98
28) bis(2-Chloroethoxy)met...	10.334	93	184960	5.190	ng	97
29) 2,4-Dichlorophenol	10.551	162	108365	4.448	ng	96
30) 1,2,4-Trichlorobenzene	10.775	180	128190	4.858	ng	99
31) Naphthalene	10.975	128	434991	5.155	ng	99
33) 4-Chloroaniline	11.098	127	186559	4.901	ng	97
34) Hexachlorobutadiene	11.210	225	74435	4.765	ng	98
35) Caprolactam	11.875	113	38690	4.345	ng	91
36) 4-Chloro-3-methylphenol	12.181	107	126057	4.655	ng	94
37) 2-Methylnaphthalene	12.575	142	309790	5.053	ng	99
38) 1-Methylnaphthalene	12.787	142	291946	5.080	ng	98
40) 1,2,4,5-Tetrachloroben...	12.916	216	146539	4.755	ng	98
41) Hexachlorocyclopentadiene	12.875	237	62826	4.059	ng	95
43) 2,4,6-Trichlorophenol	13.163	196	82195	4.156	ng	100
44) 2,4,5-Trichlorophenol	13.228	196	102205	4.260	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.575	154	396149	5.124	ng	99
47) 2-Chloronaphthalene	13.616	162	295265	5.062	ng	98
48) 2-Nitroaniline	13.839	65	69729	4.241	ng	90
49) Acenaphthylene	14.463	152	476729	4.926	ng	99
50) Dimethylphthalate	14.192	163	399188	5.051	ng	98
51) 2,6-Dinitrotoluene	14.328	165	67610	4.176	ng	# 85
52) Acenaphthene	14.798	154	288024	5.005	ng	97
53) 3-Nitroaniline	14.657	138	77327	4.224	ng	# 90
55) Dibenzofuran	15.134	168	481590	5.103	ng	98
56) 4-Nitrophenol	14.928	139	54453	3.800	ng	85
57) 2,4-Dinitrotoluene	15.104	165	77352	3.666	ng	86
58) Fluorene	15.781	166	379457	5.118	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.339	232	82945	4.203	ng	# 92
60) Diethylphthalate	15.545	149	393519	5.010	ng	98
61) 4-Chlorophenyl-phenyle...	15.775	204	185987	4.971	ng	92
62) 4-Nitroaniline	15.822	138	72659	4.000	ng	88
63) Azobenzene	16.069	77	386539	5.212	ng	93
66) n-Nitrosodiphenylamine	15.998	169	329907	5.128	ng	98
67) 4-Bromophenyl-phenylether	16.675	248	108534	4.696	ng	# 91
68) Hexachlorobenzene	16.757	284	122178	4.740	ng	92
69) Atrazine	16.945	200	94080	4.917	ng	98
71) Phenanthrene	17.539	178	605954	5.228	ng	99
72) Anthracene	17.633	178	584576	5.016	ng	98
73) Carbazole	17.904	167	549296	5.027	ng	99
74) Di-n-butylphthalate	18.427	149	599100	4.739	ng	99
75) Fluoranthene	19.492	202	650092	4.785	ng	98
77) Benzidine	19.692	184	202764	5.871	ng	99
78) Pyrene	19.851	202	687282	5.147	ng	99
80) Butylbenzylphthalate	20.716	149	229981	4.384	ng	91
81) Benzo(a)anthracene	21.568	228	670192	5.014	ng	98
82) 3,3'-Dichlorobenzidine	21.504	252	192742	4.393	ng	# 97
83) Chrysene	21.627	228	656629	5.145	ng	99
84) Bis(2-ethylhexyl)phtha...	21.463	149	354548	4.502	ng	# 96
85) Di-n-octyl phthalate	22.463	149	576684	4.354	ng	# 96
87) Indeno(1,2,3-cd)pyrene	26.956	276	688990	4.625	ng	# 95
88) Benzo(b)fluoranthene	23.374	252	622107m	4.857	ng	
89) Benzo(k)fluoranthene	23.433	252	629255	4.831	ng	97
90) Benzo(a)pyrene	24.068	252	587132	4.732	ng	98
91) Dibenzo(a,h)anthracene	26.986	278	572950	4.612	ng	# 96
92) Benzo(g,h,i)perylene	27.809	276	555765	4.642	ng	# 96

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

