

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112921\
 Data File : BP008134.D
 Acq On : 29 Nov 2021 10:26
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/30/2021
 Supervised By :Sohil Jodhani 11/30/2021

Quant Time: Nov 29 11:30:15 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 25 01:13:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	151839	20.000	ng	-0.01
21) Naphthalene-d8	10.604	136	574946	20.000	ng	0.00
39) Acenaphthene-d10	14.451	164	377117	20.000	ng	0.00
64) Phenanthrene-d10	17.210	188	793926	20.000	ng	0.00
76) Chrysene-d12	21.310	240	722658	20.000	ng	0.00
86) Perylene-d12	23.698	264	676534	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	696226	73.828	ng	-0.01
7) Phenol-d6	6.987	99	960345	75.437	ng	-0.01
23) Nitrobenzene-d5	8.969	82	964084	76.251	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	366897	76.117	ng	-0.01
45) 2-Fluorobiphenyl	13.075	172	1928984	74.298	ng	0.00
79) Terphenyl-d14	19.780	244	2867580	82.930	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.299	88	153838	34.966	ng	98
3) Pyridine	3.705	79	447470m	38.106	ng	
4) n-Nitrosodimethylamine	3.611	42	190530	38.900	ng	97
6) Aniline	7.134	93	575969	38.597	ng	98
8) 2-Chlorophenol	7.375	128	358141	37.099	ng	98
9) Benzaldehyde	6.946	77	266409	37.502	ng	96
10) Phenol	7.011	94	498226	38.085	ng	100
11) bis(2-Chloroethyl)ether	7.234	93	402494	38.493	ng	99
12) 1,3-Dichlorobenzene	7.693	146	411593	36.937	ng	99
13) 1,4-Dichlorobenzene	7.840	146	418794	37.070	ng	97
14) 1,2-Dichlorobenzene	8.158	146	397408	35.370	ng	97
15) Benzyl Alcohol	8.052	79	400589	39.716	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.334	45	611333	38.747	ng	100
17) 2-Methylphenol	8.258	107	321775	37.807	ng	99
18) Hexachloroethane	8.881	117	156150	37.172	ng	98
19) n-Nitroso-di-n-propyla...	8.622	70	327259	37.151	ng	97
20) 3+4-Methylphenols	8.587	107	449829	38.296	ng	98
22) Acetophenone	8.634	105	586767	37.311	ng	# 99
24) Nitrobenzene	9.010	77	473626	38.627	ng	98
25) Isophorone	9.540	82	844667	38.186	ng	98
26) 2-Nitrophenol	9.716	139	203926	38.547	ng	96
27) 2,4-Dimethylphenol	9.787	122	297841	37.645	ng	99
28) bis(2-Chloroethoxy)met...	10.022	93	536390	38.048	ng	99
29) 2,4-Dichlorophenol	10.257	162	361911	38.300	ng	99
30) 1,2,4-Trichlorobenzene	10.469	180	410262	37.407	ng	97
31) Naphthalene	10.651	128	1085296	36.849	ng	99
32) Benzoic acid	9.969	122	267082	41.268	ng	96
33) 4-Chloroaniline	10.769	127	449296	36.771	ng	98
34) Hexachlorobutadiene	10.940	225	285668	38.060	ng	97
35) Caprolactam	11.575	113	79907	38.069	ng	97
36) 4-Chloro-3-methylphenol	11.904	107	414117	38.337	ng	96
37) 2-Methylnaphthalene	12.269	142	770152	36.971	ng	98
38) 1-Methylnaphthalene	12.487	142	750616	37.019	ng	99
40) 1,2,4,5-Tetrachloroben...	12.640	216	474104	38.702	ng	100
41) Hexachlorocyclopentadiene	12.622	237	217669	46.217	ng	99
43) 2,4,6-Trichlorophenol	12.887	196	322967	38.913	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.957	196	343626	38.638	ng	98
46) 1,1'-Biphenyl	13.287	154	1027609	37.557	ng	99
47) 2-Chloronaphthalene	13.322	162	818371	37.874	ng	99
48) 2-Nitroaniline	13.534	65	301366	40.473	ng	97
49) Acenaphthylene	14.169	152	1251776	37.551	ng	99
50) Dimethylphthalate	13.916	163	1058420	37.375	ng	100
51) 2,6-Dinitrotoluene	14.034	165	224245	38.155	ng	100
52) Acenaphthene	14.516	154	733359	36.915	ng	98
53) 3-Nitroaniline	14.363	138	231784	38.915	ng	97
54) 2,4-Dinitrophenol	14.575	184	149354	42.981	ng	96
55) Dibenzofuran	14.851	168	1241265	36.563	ng	99
56) 4-Nitrophenol	14.687	139	186262	40.800	ng	94
57) 2,4-Dinitrotoluene	14.828	165	307464	38.120	ng	99
58) Fluorene	15.504	166	925737	33.563	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.081	232	304659m	38.523	ng	
60) Diethylphthalate	15.281	149	1051254	37.408	ng	100
61) 4-Chlorophenyl-phenyle...	15.498	204	512262	33.860	ng	100
62) 4-Nitroaniline	15.534	138	234883	38.004	ng	97
63) Azobenzene	15.792	77	1105574	38.461	ng	98
65) 4,6-Dinitro-2-methylph...	15.592	198	213530	41.120	ng	92
66) n-Nitrosodiphenylamine	15.716	169	862376	37.499	ng	100
67) 4-Bromophenyl-phenylether	16.398	248	335917	38.050	ng	99
68) Hexachlorobenzene	16.516	284	358784	38.019	ng	97
69) Atrazine	16.675	200	216466	36.267	ng	99
70) Pentachlorophenol	16.863	266	230437	41.065	ng	96
71) Phenanthrene	17.251	178	1531950	36.547	ng	100
72) Anthracene	17.345	178	1532456	36.838	ng	99
73) Carbazole	17.616	167	1474429	36.308	ng	99
74) Di-n-butylphthalate	18.175	149	1774707	34.911	ng	100
75) Fluoranthene	19.227	202	1839359	32.573	ng	99
77) Benzidine	19.404	184	449998	45.395	ng	98
78) Pyrene	19.580	202	1884921	41.193	ng	99
80) Butylbenzylphthalate	20.457	149	799904	38.871	ng	97
81) Benzo(a)anthracene	21.292	228	1791597	37.407	ng	100
82) 3,3'-Dichlorobenzidine	21.227	252	760033	33.681	ng	97
83) Chrysene	21.345	228	1691356	37.111	ng	100
84) Bis(2-ethylhexyl)phtha...	21.222	149	1046971	33.509	ng	100
85) Di-n-octyl phthalate	22.145	149	1887911	35.570	ng	99
87) Indeno(1,2,3-cd)pyrene	26.198	276	1610913	32.730	ng	99
88) Benzo(b)fluoranthene	22.974	252	1610740	39.494	ng	100
89) Benzo(k)fluoranthene	23.021	252	1537983	38.518	ng	99
90) Benzo(a)pyrene	23.598	252	1309699	37.573	ng	98
91) Dibenzo(a,h)anthracene	26.215	278	1343221	32.861	ng	98
92) Benzo(g,h,i)perylene	26.956	276	1318409	31.971	ng	98

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

