

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021624\
 Data File : BP019742.D
 Acq On : 16 Feb 2024 13:24
 Operator : MA/JU
 Sample : PB158970BSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB158970BSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/19/2024
 Supervised By :mohammad ahmed 02/19/2024

Quant Time: Feb 16 13:51:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	71684	20.000	ng	-0.01
21) Naphthalene-d8	10.704	136	248822	20.000	ng	-0.01
39) Acenaphthene-d10	14.539	164	151608	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	329252	20.000	ng	0.00
76) Chrysene-d12	21.392	240	302153	20.000	ng	0.00
86) Perylene-d12	23.815	264	359995	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	536038	120.536	ng	0.00
7) Phenol-d6	7.081	99	660830	110.206	ng	0.00
23) Nitrobenzene-d5	9.075	82	455756	79.368	ng	-0.01
42) 2,4,6-Tribromophenol	16.033	330	325519	147.055	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	995630	81.185	ng	-0.01
79) Terphenyl-d14	19.862	244	1617615	88.043	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.381	88	56814	33.159	ng	98
3) Pyridine	3.781	79	148478	31.964	ng	99
4) n-Nitrosodimethylamine	3.699	42	80946	40.025	ng	98
6) Aniline	7.234	93	137996	23.684	ng	99
8) 2-Chlorophenol	7.469	128	189717	41.552	ng	96
9) Benzaldehyde	7.051	77	10841m	2.579	ng	
10) Phenol	7.104	94	236011	39.490	ng	98
11) bis(2-Chloroethyl)ether	7.340	93	181058	38.940	ng	98
12) 1,3-Dichlorobenzene	7.787	146	230024	41.173	ng	99
13) 1,4-Dichlorobenzene	7.940	146	232771	41.125	ng	98
14) 1,2-Dichlorobenzene	8.251	146	218265	39.855	ng	99
15) Benzyl Alcohol	8.151	79	187277m	39.088	ng	
16) 2,2'-oxybis(1-Chloropr...	8.428	45	176923	38.052	ng	98
17) 2-Methylphenol	8.351	107	157945	39.255	ng	99
18) Hexachloroethane	8.969	117	87226	39.758	ng	96
19) n-Nitroso-di-n-propyla...	8.722	70	147266	35.956	ng	99
20) 3+4-Methylphenols	8.687	107	204750	37.287	ng	96
22) Acetophenone	8.734	105	293254	41.370	ng	# 98
24) Nitrobenzene	9.116	77	229671	39.724	ng	98
25) Isophorone	9.645	82	392774	39.275	ng	99
26) 2-Nitrophenol	9.822	139	101994	46.266	ng	98
27) 2,4-Dimethylphenol	9.886	122	133114	47.222	ng	98
28) bis(2-Chloroethoxy)met...	10.128	93	222699	39.359	ng	100
29) 2,4-Dichlorophenol	10.357	162	183034	42.954	ng	96
30) 1,2,4-Trichlorobenzene	10.569	180	224398	42.949	ng	100
31) Naphthalene	10.757	128	563386	41.283	ng	100
32) Benzoic acid	10.034	122	118522	43.155	ng	96
33) 4-Chloroaniline	10.881	127	49913	9.890	ng	96
34) Hexachlorobutadiene	11.022	225	168402	43.496	ng	98
35) Caprolactam	11.681	113	52107	43.195	ng	96
36) 4-Chloro-3-methylphenol	12.004	107	190558	41.013	ng	98
37) 2-Methylnaphthalene	12.363	142	383053	40.559	ng	98
38) 1-Methylnaphthalene	12.586	142	360457	38.838	ng	100
40) 1,2,4,5-Tetrachloroben...	12.727	216	269885	46.158	ng	99
41) Hexachlorocyclopentadiene	12.692	237	348299	131.863	ng	99
43) 2,4,6-Trichlorophenol	12.975	196	160448	45.318	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.051	196	171255	44.050	ng	95
46) 1,1'-Biphenyl	13.375	154	504770	41.715	ng	100
47) 2-Chloronaphthalene	13.416	162	406289	40.903	ng	100
48) 2-Nitroaniline	13.633	65	119845	43.149	ng	97
49) Acenaphthylene	14.263	152	636876	45.701	ng	100
50) Dimethylphthalate	14.010	163	489112	42.143	ng	98
51) 2,6-Dinitrotoluene	14.133	165	109663	43.006	ng	94
52) Acenaphthene	14.604	154	353686	40.893	ng	100
53) 3-Nitroaniline	14.463	138	69338	29.589	ng	97
54) 2,4-Dinitrophenol	14.674	184	142036	107.282	ng	96
55) Dibenzofuran	14.939	168	606472	43.052	ng	99
56) 4-Nitrophenol	14.774	139	172800	89.966	ng	96
57) 2,4-Dinitrotoluene	14.922	165	153544	45.950	ng	98
58) Fluorene	15.592	166	484274	43.162	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.169	232	158950	48.607	ng	96
60) Diethylphthalate	15.374	149	478001	41.407	ng	99
61) 4-Chlorophenyl-phenyle...	15.586	204	269511	42.738	ng	96
62) 4-Nitroaniline	15.627	138	103586	42.171	ng	98
63) Azobenzene	15.886	77	440587	39.468	ng	96
65) 4,6-Dinitro-2-methylph...	15.686	198	95386	50.807	ng	92
66) n-Nitrosodiphenylamine	15.810	169	416423	42.230	ng	98
67) 4-Bromophenyl-phenylether	16.486	248	178957	43.478	ng	97
68) Hexachlorobenzene	16.586	284	209786	43.645	ng	97
69) Atrazine	16.774	200	157300	48.073	ng	97
70) Pentachlorophenol	16.945	266	283415	94.683	ng	99
71) Phenanthrene	17.339	178	755957	43.628	ng	99
72) Anthracene	17.433	178	770500	44.576	ng	100
73) Carbazole	17.710	167	679607	43.280	ng	99
74) Di-n-butylphthalate	18.262	149	777463	40.948	ng	100
75) Fluoranthene	19.310	202	929972	43.934	ng	98
77) Benzidine	19.498	184	273880	43.356	ng	99
78) Pyrene	19.662	202	946589	44.389	ng	99
80) Butylbenzylphthalate	20.545	149	338600	42.519	ng	99
81) Benzo(a)anthracene	21.374	228	932254	43.831	ng	100
82) 3,3'-Dichlorobenzidine	21.309	252	289380	37.337	ng	99
83) Chrysene	21.427	228	869562	43.062	ng	99
84) Bis(2-ethylhexyl)phtha...	21.309	149	449806	37.093	ng	99
85) Di-n-octyl phthalate	22.256	149	764058	35.811	ng	99
87) Indeno(1,2,3-cd)pyrene	26.362	276	1320653	47.926	ng	97
88) Benzo(b)fluoranthene	23.074	252	954574	42.072	ng	99
89) Benzo(k)fluoranthene	23.127	252	955573	44.090	ng	99
90) Benzo(a)pyrene	23.709	252	909813	45.067	ng	98
91) Dibenzo(a,h)anthracene	26.380	278	1087080	47.965	ng	96
92) Benzo(g,h,i)perylene	27.133	276	1055457	45.904	ng	97

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

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