

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010722\
 Data File : BP008727.D
 Acq On : 07 Jan 2022 15:36
 Operator : CG/JU
 Sample : PB141824BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141824BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/10/2022
 Supervised By : Jagrut Upadhyay 01/10/2022

Quant Time: Jan 07 16:30:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 07 07:45:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	252562	20.000	ng	-0.01
21) Naphthalene-d8	10.675	136	1130813	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	960672	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	1943451	20.000	ng	0.00
76) Chrysene-d12	21.351	240	2193440	20.000	ng	0.00
86) Perylene-d12	23.780	264	2538906	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	2073770	128.580	ng	0.00
7) Phenol-d6	7.046	99	2340084	108.082	ng	0.00
23) Nitrobenzene-d5	9.028	82	1370954	68.660	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	1661268	102.500	ng	0.00
45) 2-Fluorobiphenyl	13.134	172	4276862	62.589	ng	0.00
79) Terphenyl-d14	19.822	244	7987471	74.209	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.358	88	179026	29.431	ng	98
3) Pyridine	3.752	79	421601	29.465	ng	98
4) n-Nitrosodimethylamine	3.658	42	240423	37.709	ng	91
6) Aniline	7.199	93	563348	20.712	ng	99
8) 2-Chlorophenol	7.440	128	755897	41.470	ng	99
9) Benzaldehyde	7.011	77	109488	7.472	ng	97
10) Phenol	7.069	94	905840	40.886	ng	96
11) bis(2-Chloroethyl)ether	7.293	93	628489	36.598	ng	100
12) 1,3-Dichlorobenzene	7.764	146	737606	35.819	ng	100
13) 1,4-Dichlorobenzene	7.911	146	764555	36.332	ng	99
14) 1,2-Dichlorobenzene	8.228	146	747965	36.654	ng	99
15) Benzyl Alcohol	8.111	79	610399	37.950	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.411	45	672193	34.342	ng	97
17) 2-Methylphenol	8.322	107	633726	41.490	ng	98
18) Hexachloroethane	8.958	117	252006	35.603	ng	93
19) n-Nitroso-di-n-propyla...	8.681	70	491633	34.974	ng	98
20) 3+4-Methylphenols	8.652	107	871971	41.144	ng	99
22) Acetophenone	8.693	105	1059110	35.496	ng	98
24) Nitrobenzene	9.075	77	724847	35.860	ng	100
25) Isophorone	9.605	82	1366993	34.810	ng	99
26) 2-Nitrophenol	9.787	139	405488	38.467	ng	98
27) 2,4-Dimethylphenol	9.852	122	748453	39.639	ng	99
28) bis(2-Chloroethoxy)met...	10.087	93	896685	36.621	ng	100
29) 2,4-Dichlorophenol	10.328	162	845618	41.777	ng	100
30) 1,2,4-Trichlorobenzene	10.540	180	836347	37.927	ng	99
31) Naphthalene	10.728	128	2268697	37.106	ng	100
32) Benzoic acid	10.010	122	530333	42.010	ng	98
33) 4-Chloroaniline	10.834	127	507498	18.494	ng	99
34) Hexachlorobutadiene	11.022	225	509773	38.237	ng	99
35) Caprolactam	11.622	113	270927	37.666	ng	96
36) 4-Chloro-3-methylphenol	11.963	107	839251	40.936	ng	99
37) 2-Methylnaphthalene	12.334	142	1807580	37.576	ng	99
38) 1-Methylnaphthalene	12.552	142	1674671	37.484	ng	99
40) 1,2,4,5-Tetrachloroben...	12.704	216	1053454	32.989	ng	99
41) Hexachlorocyclopentadiene	12.693	237	1229687	64.772	ng	99
43) 2,4,6-Trichlorophenol	12.946	196	745979	35.086	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.022	196	842025	36.028	ng	99
46) 1,1'-Biphenyl	13.346	154	2464862	32.914	ng	99
47) 2-Chloronaphthalene	13.387	162	1895220	32.973	ng	100
48) 2-Nitroaniline	13.587	65	473843	33.774	ng	96
49) Acenaphthylene	14.228	152	3034796	33.235	ng	100
50) Dimethylphthalate	13.969	163	2585868	33.872	ng	100
51) 2,6-Dinitrotoluene	14.081	165	580355	35.152	ng	99
52) Acenaphthene	14.575	154	2053071	37.539	ng	99
53) 3-Nitroaniline	14.410	138	365833	21.711	ng	98
54) 2,4-Dinitrophenol	14.622	184	758749	90.116	ng	99
55) Dibenzofuran	14.910	168	3457184	38.330	ng	100
56) 4-Nitrophenol	14.734	139	1177210	83.901	ng	97
57) 2,4-Dinitrotoluene	14.875	165	923196	40.967	ng	97
58) Fluorene	15.563	166	2725409	37.538	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.140	232	835115m	39.763	ng	
60) Diethylphthalate	15.340	149	2813250	38.003	ng	99
61) 4-Chlorophenyl-phenyle...	15.557	204	1457590	37.418	ng	99
62) 4-Nitroaniline	15.575	138	665065	38.014	ng	95
63) Azobenzene	15.851	77	2379445	41.006	ng	98
65) 4,6-Dinitro-2-methylph...	15.640	198	498277	46.210	ng	95
66) n-Nitrosodiphenylamine	15.769	169	2618662	44.462	ng	100
67) 4-Bromophenyl-phenylether	16.451	248	974626	42.385	ng	98
68) Hexachlorobenzene	16.569	284	1121579	41.922	ng	98
69) Atrazine	16.728	200	977438	40.885	ng	99
70) Pentachlorophenol	16.916	266	1478817	88.544	ng	100
71) Phenanthrene	17.304	178	4153086	38.462	ng	99
72) Anthracene	17.398	178	4205937	38.141	ng	99
73) Carbazole	17.669	167	3908430	38.673	ng	99
74) Di-n-butylphthalate	18.222	149	4600556	39.759	ng	99
75) Fluoranthene	19.275	202	5166538	39.957	ng	97
77) Benzidine	19.451	184	1924032	20.500	ng	98
78) Pyrene	19.622	202	5272934	36.765	ng	98
80) Butylbenzylphthalate	20.498	149	2195034	37.431	ng	99
81) Benzo(a)anthracene	21.339	228	5475434	38.050	ng	98
82) 3,3'-Dichlorobenzidine	21.269	252	1792723	27.423	ng	99
83) Chrysene	21.392	228	5303521	38.206	ng	98
84) Bis(2-ethylhexyl)phtha...	21.269	149	3306001	38.613	ng	99
85) Di-n-octyl phthalate	22.198	149	5964836	39.251	ng	98
87) Indeno(1,2,3-cd)pyrene	26.316	276	8072692	39.792	ng	97
88) Benzo(b)fluoranthene	23.039	252	6403116	38.598	ng	99
89) Benzo(k)fluoranthene	23.092	252	6168908	39.734	ng	98
90) Benzo(a)pyrene	23.674	252	6272174	39.118	ng	99
91) Dibenzo(a,h)anthracene	26.333	278	6867121	40.060	ng	98
92) Benzo(g,h,i)perylene	27.092	276	6762255	39.741	ng	98

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

