

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041322\
 Data File : BF127779.D
 Acq On : 13 Apr 2022 12:40
 Operator : CG\JU
 Sample : PB144058BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB144058BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/14/2022
 Supervised By : Jagrut Upadhyay 04/14/2022

Quant Time: Apr 13 16:49:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 16:40:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	168729	20.000	ng	0.00
21) Naphthalene-d8	8.240	136	623287	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	356446	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	631075	20.000	ng	0.00
76) Chrysene-d12	14.145	240	392712	20.000	ng	# 0.00
86) Perylene-d12	15.663	264	318653	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.593	112	756941	72.895	ng	0.01
7) Phenol-d6	6.581	99	965929	71.106	ng	0.00
23) Nitrobenzene-d5	7.528	82	983054	70.314	ng	0.00
42) 2,4,6-Tribromophenol	10.798	330	294437	77.624	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	1712693	67.974	ng	0.00
79) Terphenyl-d14	13.080	244	1843603	77.488	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.928	88	134572	27.323	ng	# 82
3) Pyridine	3.669	79	358505	27.636	ng	# 81
4) n-Nitrosodimethylamine	3.640	42	218439	31.311	ng	# 94
6) Aniline	6.622	93	453492	27.990	ng	94
8) 2-Chlorophenol	6.746	128	411739	39.089	ng	91
9) Benzaldehyde	6.510	77	295493	44.010	ng	97
10) Phenol	6.593	94	511880	37.684	ng	99
11) bis(2-Chloroethyl)ether	6.693	93	411568	36.690	ng	91
12) 1,3-Dichlorobenzene	6.898	146	430119	35.736	ng	97
13) 1,4-Dichlorobenzene	6.975	146	437974	35.998	ng	98
14) 1,2-Dichlorobenzene	7.128	146	426324	36.658	ng	98
15) Benzyl Alcohol	7.098	79	406747	34.250	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.228	45	598928	33.825	ng	96
17) 2-Methylphenol	7.204	107	340711	37.185	ng	100
18) Hexachloroethane	7.469	117	168152	36.219	ng	98
19) n-Nitroso-di-n-propyla...	7.369	70	312479	35.565	ng	96
20) 3+4-Methylphenols	7.357	107	418226	35.819	ng	# 72
22) Acetophenone	7.369	105	581622	34.726	ng	# 84
24) Nitrobenzene	7.545	77	490965	36.226	ng	96
25) Isophorone	7.781	82	911529	37.364	ng	97
26) 2-Nitrophenol	7.857	139	214619	44.065	ng	91
27) 2,4-Dimethylphenol	7.887	122	381952	40.055	ng	94
28) bis(2-Chloroethoxy)met...	7.987	93	527683	35.079	ng	99
29) 2,4-Dichlorophenol	8.098	162	372453	36.258	ng	95
30) 1,2,4-Trichlorobenzene	8.181	180	413070	34.940	ng	97
31) Naphthalene	8.263	128	1164521	35.656	ng	98
32) Benzoic acid	8.010	122	258055	37.300	ng	94
33) 4-Chloroaniline	8.310	127	167139	12.994	ng	94
34) Hexachlorobutadiene	8.375	225	275390	33.863	ng	99
35) Caprolactam	8.687	113	117550m	38.718	ng	
36) 4-Chloro-3-methylphenol	8.792	107	400888	37.224	ng	98
37) 2-Methylnaphthalene	8.957	142	790088	36.169	ng	100
38) 1-Methylnaphthalene	9.057	142	757982	35.950	ng	99
40) 1,2,4,5-Tetrachloroben...	9.122	216	416921	35.661	ng	# 97
41) Hexachlorocyclopentadiene	9.104	237	571261	81.117	ng	97
43) 2,4,6-Trichlorophenol	9.234	196	294365	37.087	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.269	196	296023	36.960	ng	99
46) 1,1'-Biphenyl	9.422	154	991728	36.380	ng	99
47) 2-Chloronaphthalene	9.451	162	773070	35.686	ng	99
48) 2-Nitroaniline	9.545	65	271655	40.375	ng	89
49) Acenaphthylene	9.869	152	1252424	38.009	ng	100
50) Dimethylphthalate	9.722	163	946591	36.637	ng	99
51) 2,6-Dinitrotoluene	9.787	165	206767	42.456	ng	90
52) Acenaphthene	10.039	154	838853	40.408	ng	97
53) 3-Nitroaniline	9.957	138	106425	20.685	ng	91
54) 2,4-Dinitrophenol	10.063	184	215158	88.529	ng	97
55) Dibenzofuran	10.210	168	1074096	35.456	ng	97
56) 4-Nitrophenol	10.116	139	304127	75.257	ng	88
57) 2,4-Dinitrotoluene	10.192	165	276584	45.807	ng	# 83
58) Fluorene	10.557	166	822666	36.753	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	259490	37.181	ng	100
60) Diethylphthalate	10.422	149	890771	36.028	ng	99
61) 4-Chlorophenyl-phenyle...	10.545	204	437542	35.576	ng	99
62) 4-Nitroaniline	10.575	138	189746	39.158	ng	95
63) Azobenzene	10.704	77	936610	33.901	ng	96
65) 4,6-Dinitro-2-methylph...	10.598	198	138364	45.393	ng	87
66) n-Nitrosodiphenylamine	10.663	169	749665	37.718	ng	99
67) 4-Bromophenyl-phenylether	11.033	248	289096	37.746	ng	# 91
68) Hexachlorobenzene	11.098	284	321558	38.407	ng	95
69) Atrazine	11.186	200	277405	42.008	ng	97
70) Pentachlorophenol	11.292	266	352325	70.459	ng	98
71) Phenanthrene	11.522	178	1213344	36.288	ng	99
72) Anthracene	11.575	178	1246663	37.692	ng	98
73) Carbazole	11.728	167	1092484	35.251	ng	99
74) Di-n-butylphthalate	12.051	149	1327679	35.837	ng	99
75) Fluoranthene	12.710	202	1325806	36.337	ng	98
77) Benzidine	12.833	184	497276	48.514	ng	98
78) Pyrene	12.945	202	1329107	40.079	ng	99
80) Butylbenzylphthalate	13.551	149	474202	38.405	ng	96
81) Benzo(a)anthracene	14.133	228	1014248	38.669	ng	98
82) 3,3'-Dichlorobenzidine	14.092	252	181571	22.205	ng	# 98
83) Chrysene	14.169	228	915499	36.362	ng	99
84) Bis(2-ethylhexyl)phtha...	14.110	149	603160	34.656	ng	# 97
85) Di-n-octyl phthalate	14.727	149	826293	32.470	ng	# 93
87) Indeno(1,2,3-cd)pyrene	17.221	276	801519	40.259	ng	99
88) Benzo(b)fluoranthene	15.210	252	805713	39.305	ng	99
89) Benzo(k)fluoranthene	15.239	252	791157	38.914	ng	100
90) Benzo(a)pyrene	15.598	252	757188	45.446	ng	99
91) Dibenzo(a,h)anthracene	17.245	278	696477	43.176	ng	98
92) Benzo(g,h,i)perylene	17.704	276	700378	42.351	ng	98

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

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