

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111722\
 Data File : BF131279.D
 Acq On : 17 Nov 2022 11:41
 Operator : CG\JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

11/18/2022

Supervised By :Jagrut
 Upadhyay

11/18/2022

Quant Time: Nov 18 00:54:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 18 00:51:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	89179	20.000	ng	0.00
21) Naphthalene-d8	8.257	136	329638	20.000	ng	0.00
39) Acenaphthene-d10	10.027	164	186894	20.000	ng	0.00
64) Phenanthrene-d10	11.521	188	344929	20.000	ng	0.00
76) Chrysene-d12	14.174	240	239789	20.000	ng	0.00
86) Perylene-d12	15.715	264	171673	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	49011	9.732	ng	-0.01
7) Phenol-d6	6.569	99	64152	10.057	ng	-0.02
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	10.810	330	12161	6.812	ng	-0.01
45) 2-Fluorobiphenyl	9.339	172	143817	11.261	ng	0.00
79) Terphenyl-d14	13.110	244	154047	11.032	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.787	88	12036	5.216	ng	99
3) Pyridine	3.557	79	32402	4.968	ng	96
4) n-Nitrosodimethylamine	3.498	42	19791	5.206	ng	94
6) Aniline	6.622	93	40222	5.027	ng	99
8) 2-Chlorophenol	6.745	128	26259	4.638	ng	97
9) Benzaldehyde	6.516	77	26093	5.639	ng	100
10) Phenol	6.586	94	34897	4.890	ng	93
11) bis(2-Chloroethyl)ether	6.692	93	29557	5.290	ng	99
12) 1,3-Dichlorobenzene	6.910	146	34608m	5.404	ng	
13) 1,4-Dichlorobenzene	6.986	146	35232	5.439	ng	97
14) 1,2-Dichlorobenzene	7.139	146	33596	5.612	ng	99
15) Benzyl Alcohol	7.098	79	25582	4.617	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.239	45	55486	5.272	ng	99
17) 2-Methylphenol	7.204	107	25023	5.177	ng	89
18) Hexachloroethane	7.486	117	10986	4.665	ng	94
19) n-Nitroso-di-n-propyla...	7.369	70	23376	5.259	ng	97
20) 3+4-Methylphenols	7.357	107	31322	5.010	ng	92
22) Acetophenone	7.375	105	45454	5.290	ng	# 98
24) Nitrobenzene	7.545	77	22301	3.518	ng	98
25) Isophorone	7.786	82	56890	4.849	ng	98
26) 2-Nitrophenol	7.869	139	4935	2.278	ng	95
27) 2,4-Dimethylphenol	7.898	122	21130	4.400	ng	100
28) bis(2-Chloroethoxy)met...	7.998	93	33735	5.100	ng	98
29) 2,4-Dichlorophenol	8.104	162	20520	3.951	ng	97
30) 1,2,4-Trichlorobenzene	8.198	180	30980	5.088	ng	97
31) Naphthalene	8.280	128	95789	5.364	ng	99
33) 4-Chloroaniline	8.328	127	34565	4.917	ng	96
34) Hexachlorobutadiene	8.398	225	20324	5.274	ng	97
35) Caprolactam	8.651	113	5361	3.493	ng	90
36) 4-Chloro-3-methylphenol	8.792	107	23436	4.287	ng	98
37) 2-Methylnaphthalene	8.975	142	62138	5.224	ng	94
38) 1-Methylnaphthalene	9.075	142	60919	5.282	ng	98
40) 1,2,4,5-Tetrachloroben...	9.139	216	32095	5.334	ng	99
43) 2,4,6-Trichlorophenol	9.245	196	13863	3.524	ng	99
44) 2,4,5-Trichlorophenol	9.280	196	15150	3.589	ng	95
46) 1,1'-Biphenyl	9.439	154	76418	5.319	ng	98

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47) 2-Chloronaphthalene	9.463	162	60706	5.216	ng	98
48) 2-Nitroaniline	9.557	65	8711	2.711	ng	98
49) Acenaphthylene	9.886	152	92033	5.178	ng	99
50) Dimethylphthalate	9.733	163	67522	5.045	ng	99
51) 2,6-Dinitrotoluene	9.798	165	6321	2.710	ng	91
52) Acenaphthene	10.057	154	56112	5.058	ng	99
53) 3-Nitroaniline	9.969	138	7017	2.785	ng	# 87
55) Dibenzofuran	10.227	168	84636	5.325	ng	98
57) 2,4-Dinitrotoluene	10.198	165	6543	2.364	ng	97
58) Fluorene	10.574	166	67304	5.381	ng	96
59) 2,3,4,6-Tetrachlorophenol	10.339	232	12521	3.501	ng	99
60) Diethylphthalate	10.439	149	62017	4.902	ng	97
61) 4-Chlorophenyl-phenyle...	10.569	204	35062	5.669	ng	91
62) 4-Nitroaniline	10.574	138	8014	3.220	ng	82
63) Azobenzene	10.727	77	68071	5.139	ng	97
66) n-Nitrosodiphenylamine	10.680	169	56804	5.183	ng	95
67) 4-Bromophenyl-phenylether	11.063	248	20353	5.026	ng	91
68) Hexachlorobenzene	11.121	284	22782	5.191	ng	96
69) Atrazine	11.204	200	16666	4.640	ng	100
71) Phenanthrene	11.545	178	100540	5.282	ng	99
72) Anthracene	11.598	178	96563	5.193	ng	99
73) Carbazole	11.745	167	81338	4.886	ng	98
74) Di-n-butylphthalate	12.080	149	83000	4.284	ng	99
75) Fluoranthene	12.739	202	104895	4.996	ng	99
77) Benzidine	12.863	184	32141	6.255	ng	98
78) Pyrene	12.968	202	108894	5.268	ng	99
80) Butylbenzylphthalate	13.592	149	23868	3.407	ng	93
81) Benzo(a)anthracene	14.162	228	87899	5.275	ng	99
82) 3,3'-Dichlorobenzidine	14.127	252	19072	3.911	ng	# 99
83) Chrysene	14.204	228	82685	5.093	ng	98
84) Bis(2-ethylhexyl)phtha...	14.157	149	27989	3.177	ng	# 99
87) Indeno(1,2,3-cd)pyrene	17.286	276	46144	4.069	ng	98
88) Benzo(b)fluoranthene	15.251	252	56804	4.739	ng	100
89) Benzo(k)fluoranthene	15.286	252	61822	5.203	ng	99
90) Benzo(a)pyrene	15.645	252	43931	4.585	ng	98
91) Dibenzo(a,h)anthracene	17.315	278	38055	4.090	ng	97
92) Benzo(g,h,i)perylene	17.768	276	38183	4.112	ng	92

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

