

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121821\
 Data File : BF126257.D
 Acq On : 18 Dec 2021 16:42
 Operator : JU/CG
 Sample : M5082-02MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2-121-1(55-62)MSD

Quant Time: Dec 20 08:03:02 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 12/20/2021
 Supervised By :mohammad ahmed 12/29/2021

12/20/2021
 Supervised By :mohammad
 ahmed

12/29/2021

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	229758	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	1010066	20.000	ng	# 0.00
39) Acenaphthene-d10	9.857	164	471115	20.000	ng	0.00
64) Phenanthrene-d10	11.339	188	753292	20.000	ng	# 0.00
76) Chrysene-d12	13.980	240	393768	20.000	ng	# 0.01
86) Perylene-d12	15.421	264	374210	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1577380	101.103	ng	0.02
7) Phenol-d6	6.451	99	2086378	105.319	ng	0.00
23) Nitrobenzene-d5	7.381	82	1363180	73.093	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	410416	108.401	ng	0.01
45) 2-Fluorobiphenyl	9.180	172	1778047	66.194	ng	0.00
79) Terphenyl-d14	12.927	244	1524178	67.278	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.657	88	332527	43.663	ng	# 100
3) Pyridine	3.381	79	761458	37.650	ng	# 79
4) n-Nitrosodimethylamine	3.340	42	383705	49.364	ng	# 1
6) Aniline	6.481	93	1079546	42.902	ng	96
8) 2-Chlorophenol	6.604	128	809186	51.145	ng	90
9) Benzaldehyde	6.363	77	587150	49.117	ng	95
10) Phenol	6.463	94	1074798	48.334	ng	94
11) bis(2-Chloroethyl)ether	6.557	93	832888	48.253	ng	98
12) 1,3-Dichlorobenzene	6.757	146	723527	45.335	ng	96
13) 1,4-Dichlorobenzene	6.834	146	727608	45.146	ng	95
14) 1,2-Dichlorobenzene	6.986	146	718515	48.217	ng	95
15) Benzyl Alcohol	6.963	79	725654	50.177	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.098	45	1370496	48.519	ng	94
17) 2-Methylphenol	7.075	107	690633	49.829	ng	98
18) Hexachloroethane	7.334	117	300664	47.427	ng	95
19) n-Nitroso-di-n-propyla...	7.239	70	565088	46.466	ng	# 71
20) 3+4-Methylphenols	7.228	107	729401	44.515	ng	# 84
22) Acetophenone	7.234	105	1028970	44.264	ng	96
24) Nitrobenzene	7.398	77	884248	47.495	ng	90
25) Isophorone	7.645	82	1675371	47.571	ng	# 88
26) 2-Nitrophenol	7.716	139	426622	50.831	ng	# 82
27) 2,4-Dimethylphenol	7.757	122	764651	53.648	ng	97
28) bis(2-Chloroethoxy)met...	7.851	93	1055619	45.589	ng	# 97
29) 2,4-Dichlorophenol	7.957	162	596606	47.090	ng	92
30) 1,2,4-Trichlorobenzene	8.045	180	583998	44.563	ng	99
31) Naphthalene	8.122	128	2161286	45.755	ng	99
32) Benzoic acid	7.892	122	587522	47.847	ng	95
33) 4-Chloroaniline	8.169	127	552243	28.000	ng	96
34) Hexachlorobutadiene	8.245	225	277496	43.115	ng	99
35) Caprolactam	8.551	113	239202m	75.969	ng	
36) 4-Chloro-3-methylphenol	8.657	107	715165	47.853	ng	89
37) 2-Methylnaphthalene	8.816	142	1374813	47.261	ng	99
38) 1-Methylnaphthalene	8.916	142	1269443	44.967	ng	96
40) 1,2,4,5-Tetrachloroben...	8.986	216	436419	42.921	ng	# 96
41) Hexachlorocyclopentadiene	8.975	237	540774	92.874	ng	93
43) 2,4,6-Trichlorophenol	9.092	196	370420	46.311	ng	94

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44) 2,4,5-Trichlorophenol	9.133	196	372442	44.964	ng	94
46) 1,1'-Biphenyl	9.280	154	1566034	44.956	ng	96
47) 2-Chloronaphthalene	9.304	162	1170630	44.665	ng	97
48) 2-Nitroaniline	9.398	65	471064	49.242	ng	89
49) Acenaphthylene	9.722	152	1821461	45.482	ng	97
50) Dimethylphthalate	9.586	163	1355747	45.429	ng	98
51) 2,6-Dinitrotoluene	9.645	165	312322	48.352	ng	# 87
52) Acenaphthene	9.892	154	1286155	49.522	ng	100
53) 3-Nitroaniline	9.810	138	270891	32.995	ng	# 75
54) 2,4-Dinitrophenol	9.916	184	309232	116.269	ng	# 82
55) Dibenzofuran	10.069	168	1523682	43.090	ng	98
56) 4-Nitrophenol	9.975	139	608619	102.629	ng	# 78
57) 2,4-Dinitrotoluene	10.051	165	417626	50.685	ng	# 96
58) Fluorene	10.410	166	1081016	42.229	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.180	232	283190	46.110	ng	# 99
60) Diethylphthalate	10.286	149	1333658	44.469	ng	98
61) 4-Chlorophenyl-phenyle...	10.398	204	475594	41.263	ng	98
62) 4-Nitroaniline	10.427	138	375801	54.566	ng	# 73
63) Azobenzene	10.563	77	1639883	45.144	ng	84
65) 4,6-Dinitro-2-methylph...	10.457	198	196706	47.820	ng	89
66) n-Nitrosodiphenylamine	10.522	169	1046752	43.298	ng	96
67) 4-Bromophenyl-phenylether	10.892	248	314643	44.505	ng	97
68) Hexachlorobenzene	10.957	284	370630	45.582	ng	# 85
69) Atrazine	11.051	200	324621	73.271	ng	96
70) Pentachlorophenol	11.151	266	384012	85.130	ng	91
71) Phenanthrene	11.369	178	1687829	43.411	ng	97
72) Anthracene	11.422	178	1713267	43.903	ng	98
73) Carbazole	11.574	167	1595966	43.427	ng	100
74) Di-n-butylphthalate	11.910	149	1992890	42.776	ng	100
75) Fluoranthene	12.557	202	1580834	45.128	ng	93
77) Benzidine	12.674	184	854898	137.710	ng	98
78) Pyrene	12.780	202	1596092	44.120	ng	96
80) Butylbenzylphthalate	13.404	149	801477	47.080	ng	# 80
81) Benzo(a)anthracene	13.968	228	977756	41.919	ng	100
82) 3,3'-Dichlorobenzidine	13.927	252	310277	29.118	ng	# 96
83) Chrysene	14.004	228	1108930	45.759	ng	98
84) Bis(2-ethylhexyl)phtha...	13.968	149	875557	46.025	ng	100
85) Di-n-octyl phthalate	14.574	149	1683021	49.629	ng	99
87) Indeno(1,2,3-cd)pyrene	16.856	276	1135907	45.509	ng	# 92
88) Benzo(b)fluoranthene	15.004	252	1158634m	51.530	ng	
89) Benzo(k)fluoranthene	15.039	252	977600	45.450	ng	# 96
90) Benzo(a)pyrene	15.362	252	1033709	55.316	ng	# 95
91) Dibenzo(a,h)anthracene	16.880	278	956663	47.149	ng	96
92) Benzo(g,h,i)perylene	17.292	276	973379	46.366	ng	92

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

