

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
 Data File : BF136989.D
 Acq On : 08 Jan 2024 10:39
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 01/09/2024

Supervised By :mohammad ahmed 01/09/2024

Quant Time: Jan 08 11:00:50 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Dec 21 01:54:25 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	71646	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	273505	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	138981	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	242393	20.000	ng	0.00
76) Chrysene-d12	13.974	240	113262	20.000	ng	# 0.00
86) Perylene-d12	15.457	264	123957	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	369855	80.546	ng	0.00
7) Phenol-d6	6.445	99	459521	77.231	ng	0.00
23) Nitrobenzene-d5	7.357	82	421273	78.818	ng	0.00
42) 2,4,6-Tribromophenol	10.622	330	120481	73.607	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	832896	80.603	ng	0.00
79) Terphenyl-d14	12.910	244	709412	87.530	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.534	88	75479	39.451	ng	94
3) Pyridine	3.299	79	179734	38.503	ng	91
4) n-Nitrosodimethylamine	3.275	42	90807	36.427	ng	97
6) Aniline	6.457	93	217073	36.490	ng	# 29
8) 2-Chlorophenol	6.581	128	201286	40.057	ng	94
9) Benzaldehyde	6.345	77	126618	37.413	ng	98
10) Phenol	6.457	94	240604	37.881	ng	91
11) bis(2-Chloroethyl)ether	6.528	93	187561	38.830	ng	99
12) 1,3-Dichlorobenzene	6.728	146	217385	39.656	ng	97
13) 1,4-Dichlorobenzene	6.804	146	221549	39.554	ng	99
14) 1,2-Dichlorobenzene	6.957	146	205356	39.401	ng	99
15) Benzyl Alcohol	6.939	79	155455	38.728	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.063	45	290505	36.786	ng	71
17) 2-Methylphenol	7.057	107	155287	37.303	ng	# 89
18) Hexachloroethane	7.292	117	78561	38.621	ng	90
19) n-Nitroso-di-n-propyla...	7.204	70	139249	38.607	ng	98
20) 3+4-Methylphenols	7.210	107	203483	39.126	ng	# 54
22) Acetophenone	7.198	105	276962	40.401	ng	# 88
24) Nitrobenzene	7.375	77	208964	39.028	ng	96
25) Isophorone	7.610	82	375292	38.588	ng	98
26) 2-Nitrophenol	7.686	139	105948	41.359	ng	97
27) 2,4-Dimethylphenol	7.728	122	127638	40.927	ng	98
28) bis(2-Chloroethoxy)met...	7.816	93	233622	39.518	ng	100
29) 2,4-Dichlorophenol	7.934	162	162614	40.501	ng	95
30) 1,2,4-Trichlorobenzene	8.004	180	182888	40.882	ng	97
31) Naphthalene	8.086	128	598684	39.844	ng	99
32) Benzoic acid	7.875	122	115842m	38.389	ng	
33) 4-Chloroaniline	8.139	127	214184	38.607	ng	97
34) Hexachlorobutadiene	8.198	225	110947	40.820	ng	98
35) Caprolactam	8.528	113	55181m	41.660	ng	
36) 4-Chloro-3-methylphenol	8.633	107	175664	38.863	ng	96
37) 2-Methylnaphthalene	8.781	142	385179	39.832	ng	99
38) 1-Methylnaphthalene	8.875	142	375351	39.564	ng	100
40) 1,2,4,5-Tetrachloroben...	8.945	216	179900	41.797	ng	99
41) Hexachlorocyclopentadiene	8.928	237	59949	45.907	ng	99
43) 2,4,6-Trichlorophenol	9.063	196	110570	40.057	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
 Data File : BF136989.D
 Acq On : 08 Jan 2024 10:39
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 08 11:00:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 01/09/2024
 Supervised By :mohammad ahmed 01/09/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.116	196	119691	38.911	ng	99
46) 1,1'-Biphenyl	9.245	154	471478	40.426	ng	99
47) 2-Chloronaphthalene	9.269	162	359787	40.580	ng	99
48) 2-Nitroaniline	9.375	65	104033	37.470	ng	92
49) Acenaphthylene	9.686	152	529107	39.043	ng	99
50) Dimethylphthalate	9.545	163	387205	38.143	ng	98
51) 2,6-Dinitrotoluene	9.616	165	89037	38.646	ng	89
52) Acenaphthene	9.857	154	332255	39.552	ng	99
53) 3-Nitroaniline	9.792	138	91118	37.335	ng	87
54) 2,4-Dinitrophenol	9.904	184	44673	37.088	ng	90
55) Dibenzofuran	10.033	168	486868	38.234	ng	96
56) 4-Nitrophenol	9.969	139	54445	33.047	ng	98
57) 2,4-Dinitrotoluene	10.028	165	108593	36.308	ng	# 95
58) Fluorene	10.375	166	392447	38.841	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.157	232	83055	35.076	ng	96
60) Diethylphthalate	10.251	149	389043	36.936	ng	97
61) 4-Chlorophenyl-phenyle...	10.363	204	190665	38.823	ng	99
62) 4-Nitroaniline	10.410	138	81276m	34.545	ng	
63) Azobenzene	10.527	77	359374	36.530	ng	98
65) 4,6-Dinitro-2-methylph...	10.439	198	53371	38.420	ng	94
66) n-Nitrosodiphenylamine	10.492	169	332977	42.288	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	114459	42.519	ng	96
68) Hexachlorobenzene	10.927	284	127015	42.339	ng	93
69) Atrazine	11.022	200	77665	38.564	ng	95
70) Pentachlorophenol	11.127	266	51029	36.493	ng	99
71) Phenanthrene	11.339	178	497412	39.993	ng	99
72) Anthracene	11.392	178	504119	40.023	ng	99
73) Carbazole	11.557	167	451120	37.988	ng	99
74) Di-n-butylphthalate	11.880	149	565485	39.023	ng	99
75) Fluoranthene	12.533	202	492139	36.993	ng	97
77) Benzidine	12.663	184	161879	48.472	ng	99
78) Pyrene	12.769	202	482871	43.429	ng	100
80) Butylbenzylphthalate	13.386	149	169049	41.785	ng	99
81) Benzo(a)anthracene	13.963	228	319927	40.678	ng	98
82) 3,3'-Dichlorobenzidine	13.927	252	94651	42.377	ng	98
83) Chrysene	13.998	228	307595	39.994	ng	99
84) Bis(2-ethylhexyl)phtha...	13.945	149	222654	43.741	ng	98
85) Di-n-octyl phthalate	14.563	149	327154	41.549	ng	# 100
87) Indeno(1,2,3-cd)pyrene	16.986	276	406720	45.444	ng	100
88) Benzo(b)fluoranthene	15.021	252	313215	37.832	ng	98
89) Benzo(k)fluoranthene	15.051	252	270397	34.564	ng	99
90) Benzo(a)pyrene	15.398	252	273227	39.097	ng	100
91) Dibenzo(a,h)anthracene	16.998	278	332480	45.282	ng	97
92) Benzo(g,h,i)perylene	17.445	276	344431	45.800	ng	95

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

